

**AMENDMENT**

Subject matter to be added is in bold and underlined.

Subject matter to be deleted is in bold and strikethrough.

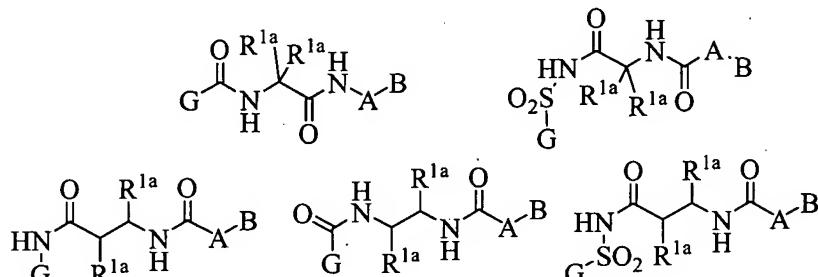
**In the Claims:**

Please enter rewritten Claims 1-5 and 7-9 and new Claims 15-22 as follows.

This listing of claims will replace all prior versions and listings of claims in the application.

**Listing of Claims:**

1. (Currently Amended) A compound of selected from: **formula I:**



**P-M-M<sub>1</sub>**

**I**

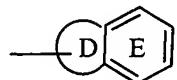
or a stereoisomer or pharmaceutically acceptable salt thereof, wherein;

~~one of P and M<sub>1</sub> is G and the other A-B;~~

G is a group of formula IIa or IIb:



IIa



IIb

ring D, including the two atoms of Ring E to which it is attached, is a 5-6 membered ring consisting of: carbon atoms and 0-3 heteroatoms selected from the group consisting of N, O, and S(O)p;

ring D is substituted with 0-2 R, 0-2 carbonyls, and there are 0-3 ring double bonds;

E is selected from phenyl, pyridyl, pyrimidyl, pyrazinyl, and pyridazinyl, and is substituted with 1-2 R;

alternatively, ring D is absent and ring E is selected from phenyl, pyridyl, pyrimidyl, pyrazinyl, pyridazinyl, pyrrolyl, pyrazolyl, imidazolyl, isoxazolyl, oxazolyl, triazolyl, thienyl, and thiazolyl, and ring E is substituted with 1-2 R;

alternatively, ring D is absent and ring E is selected from phenyl, pyridyl, pyrimidyl, pyrazinyl, pyridazinyl, pyrrolyl, pyrazolyl, imidazolyl, isoxazolyl, oxazolyl, triazolyl, thienyl, and thiazolyl, and ring E is substituted with 1 R and with a 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)p, wherein the 5-6 membered heterocycle is substituted with 0-1 carbonyls and 1-2 R and there are 0-3 ring double bonds;

R is selected from H, C<sub>1-4</sub> alkyl, F, Cl, Br, I, OH, OCH<sub>3</sub>, OCH<sub>2</sub>CH<sub>3</sub>, OCH(CH<sub>3</sub>)<sub>2</sub>, QCH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, -CN, C(=NR<sup>8</sup>)NR<sup>7</sup>R<sup>9</sup>, NHC(=NR<sup>8</sup>)NR<sup>7</sup>R<sup>9</sup>, ONHC(=NR<sup>8</sup>)NR<sup>7</sup>R<sup>9</sup>, NR<sup>8</sup>CH(=NR<sup>7</sup>), NH<sub>2</sub>, NH(C<sub>1-3</sub> alkyl), N(C<sub>1-3</sub> alkyl)<sub>2</sub>, C(=NH)NH<sub>2</sub>, CH<sub>2</sub>NH<sub>2</sub>, CH<sub>2</sub>NH(C<sub>1-3</sub> alkyl), CH<sub>2</sub>N(C<sub>1-3</sub> alkyl)<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>NH<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>NH(C<sub>1-3</sub> alkyl), CH<sub>2</sub>CH<sub>2</sub>N(C<sub>1-3</sub> alkyl)<sub>2</sub>, (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>C(O)H, (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>C(O)R<sup>2c</sup>, (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>NR<sup>7</sup>R<sup>8</sup>, (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>C(O)NR<sup>7</sup>R<sup>8</sup>, (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>NR<sup>7</sup>C(O)R<sup>7</sup>, (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>OR<sup>3</sup>, (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>S(O)pNR<sup>7</sup>R<sup>8</sup>, (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>NR<sup>7</sup>S(O)pR<sup>7</sup>, (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>SR<sup>3</sup>, (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>S(O)R<sup>3</sup>, (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>S(O)R<sup>3</sup>, and OCF<sub>3</sub>;

alternatively, when 2 R groups are attached to adjacent atoms, they combine to form methylenedioxy or ethylenedioxy;

~~M is 3-8 membered linear chain consisting of: carbon atoms, 0-3 carbonyl groups, 0-1 thiocarbonyl groups, and 1-3 heteroatoms selected from O, N, and S(O)p; and M is substituted with 0-3 R<sup>1a</sup> and 0-2 R<sup>2</sup>, and there are 0-2 double bonds and 0-1 triple bond; provided that other than an S-S, S-O, or O-O bond is present in M;~~

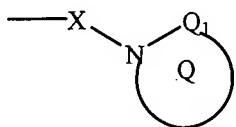
~~provided that linker M comprises other than a N-C(O)-C(O)-N group;~~

~~further provided that one or more of the following apply:~~

- (a) if linker M comprises a ureido-methylene-carbonyl-amino or carbamoyloxy-methylene-carbonyl-amino group, then ring D is present or ring E is other than phenyl or pyridyl;
- (b) there is at least one S(O)<sub>p</sub> group present in linker M;
- (c) there are at least two carbonyl groups present in linker M;
- (d) ring D is present in group G;
- (e) ring E is other than phenyl; and
- (f) if ring D is absent and ring E is phenyl, then R is other than CN, C(=NR<sup>8</sup>)NR<sup>7</sup>R<sup>9</sup>, NR<sup>8</sup>CH(=NR<sup>7</sup>), NH<sub>2</sub>, NH(C<sub>1-3</sub>-alkyl), N(C<sub>1-3</sub>-alkyl)<sub>2</sub>, C(=NH)NH<sub>2</sub>, CH<sub>2</sub>NH<sub>2</sub>, CH<sub>2</sub>NH(C<sub>1-3</sub>-alkyl), CH<sub>2</sub>N(C<sub>1-3</sub>-alkyl)<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>NH<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>NH(C<sub>1-3</sub>-alkyl), CH<sub>2</sub>CH<sub>2</sub>N(C<sub>1-3</sub>-alkyl)<sub>2</sub>, (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>NR<sup>7</sup>R<sup>8</sup>, and (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>C(O)NR<sup>7</sup>R<sup>8</sup>;

A is selected from:

C<sub>3-10</sub> carbocycle substituted with 0-2 R<sup>4</sup>, and  
5-12 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)p and substituted with 0-2 R<sup>4</sup>;



B is ; provided that Z and B are attached to different atoms on A and that the A-X-N moiety forms other than a N-N-N group;

Q<sub>1</sub> is selected from C=O and SO<sub>2</sub>;

ring Q is a 4-8 membered monocyclic or bicyclic ring consisting of, in addition to the N-Q<sub>1</sub> group shown, carbon atoms and 0-2 heteroatoms selected from NR<sup>4c</sup>, O, S, S(O), and S(O)<sub>2</sub>, wherein:

0-2 double bonds are present within the ring and the ring is substituted with 0-2 R<sup>4a</sup>;

alternatively, ring Q is a 4-8 membered monocyclic or bicyclic ring to which another ring is fused, wherein:

the 4-7 membered ring consists of, in addition to the shown amide group, carbon atoms and 0-2 heteroatoms selected from NR<sup>4c</sup>, O, S, S(O), and S(O)<sub>2</sub> and 0-2 double bonds are present within the ring;

the fusion ring is phenyl or a 5-6 membered heteroaromatic consisting of carbon atoms and 1-2 heteroatoms selected from NR<sup>4c</sup>, O, S, S(O), and S(O)<sub>2</sub>;

ring Q, which includes the 4-7 membered ring and the fusion ring, is substituted with 0-3 R<sup>4a</sup>;

alternatively, two non-adjacent atoms of one of the rings of ring Q are bridged with 1-2 atoms selected from: carbon atoms, NR<sup>4c</sup>, O, S, S(O), and S(O)<sub>2</sub>, provided bonds other than O-O, S(O)<sub>p</sub>-O, S(O)<sub>p</sub>-S(O)<sub>p</sub>, N-O, and N-S(O)<sub>p</sub> are present;

X is absent or is selected from -(CR<sup>2</sup>R<sup>2a</sup>)<sub>1-4</sub>-, -CR<sup>2</sup>(CR<sup>2</sup>R<sup>2b</sup>)(CH<sub>2</sub>)<sub>t</sub>-, -C(O)-, -C(=NR<sup>1c</sup>)-, -CR<sup>2</sup>(NR<sup>1c</sup>R<sup>2</sup>)-, -CR<sup>2</sup>(OR<sup>2</sup>)-, -CR<sup>2</sup>(SR<sup>2</sup>)-, -C(O)CR<sup>2</sup>R<sup>2a</sup>-,

-CR<sup>2</sup>R<sup>2a</sup>C(O), -S(O)-, -S(O)<sub>2</sub>-, -SCR<sup>2</sup>R<sup>2a</sup>-, -S(O)CR<sup>2</sup>R<sup>2a</sup>-, -S(O)<sub>2</sub>CR<sup>2</sup>R<sup>2a</sup>-,  
-CR<sup>2</sup>R<sup>2a</sup>S(O)-, -CR<sup>2</sup>R<sup>2a</sup>S(O)<sub>2</sub>-, -S(O)<sub>2</sub>NR<sup>2</sup>CR<sup>2</sup>R<sup>2a</sup>-, --NR<sup>2</sup>S(O)<sub>2</sub>-,  
-CR<sup>2</sup>R<sup>2a</sup>NR<sup>2</sup>S(O)<sub>2</sub>-, -NR<sup>2</sup>S(O)<sub>2</sub>CR<sup>2</sup>R<sup>2a</sup>-, -NR<sup>2</sup>C(O)-, -C(O)NR<sup>2</sup>CR<sup>2</sup>R<sup>2a</sup>-,  
-NR<sup>2</sup>C(O)CR<sup>2</sup>R<sup>2a</sup>-, -CR<sup>2</sup>R<sup>2a</sup>NR<sup>2</sup>C(O)-, -NR<sup>2</sup>CR<sup>2</sup>R<sup>2a</sup>-, and -OCR<sup>2</sup>R<sup>2a</sup>;

R<sup>1a</sup>, at each occurrence, is selected from H, -(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1b</sup>, -(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>CR<sup>3</sup>R<sup>1b</sup>R<sup>1b</sup>,  
-(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>O-(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1b</sup>, -C<sub>2-6</sub> alkenylene-R<sup>1b</sup>, -C<sub>2-6</sub> alkynylene-R<sup>1b</sup>,  
-(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>C(=NR<sup>1b</sup>)NR<sup>3</sup>R<sup>1b</sup>, NR<sup>3</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1c</sup>, O(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1c</sup>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>SCR<sup>3</sup>R<sup>3a</sup>R<sup>1c</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>NR<sup>3</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1b</sup>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>C(O)NR<sup>2</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1b</sup>, CO<sub>2</sub>(CR<sup>3</sup>R<sup>3a</sup>)<sub>t</sub>R<sup>1b</sup>, O(CR<sup>3</sup>R<sup>3a</sup>)<sub>t</sub>R<sup>1b</sup>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>S(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1b</sup>, S(O)p(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>, O(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>,  
NR<sup>3</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>, OC(O)NR<sup>3</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>, NR<sup>3</sup>C(O)NR<sup>3</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>,  
NR<sup>3</sup>C(O)O(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>, and NR<sup>3</sup>C(O)(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>, provided that R<sup>1a</sup> forms  
other than an N-halo, N-S, O-O, or N-CN bond;

alternatively, when two R<sup>1a</sup> groups are attached to the same carbon atom, together with  
the carbon atom to which they are attached they form a 3-10 membered  
carbocyclic or heterocyclic ring consisting of: carbon atoms and 0-4 heteroatoms  
selected from the group consisting of N, O, and S(O)<sub>p</sub>, this ring being substituted  
with 0-2 R<sup>4</sup> and 0-3 ring double bonds;

R<sup>1b</sup> is selected from H, C<sub>1-3</sub> alkyl, F, Cl, Br, I, -CN, -NO<sub>2</sub>, -CHO, (CF<sub>2</sub>)<sub>r</sub>CF<sub>3</sub>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>OR<sup>2</sup>, NR<sup>2</sup>R<sup>2a</sup>, C(O)R<sup>2b</sup>, CO<sub>2</sub>R<sup>2b</sup>, OC(O)R<sup>2</sup>, (CF<sub>2</sub>)<sub>r</sub>CO<sub>2</sub>R<sup>2a</sup>,  
S(O)<sub>p</sub>R<sup>2b</sup>, NR<sup>2</sup>(CH<sub>2</sub>)<sub>r</sub>OR<sup>2</sup>, C(=NR<sup>2c</sup>)NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>,  
NR<sup>2</sup>C(O)NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>C(O)<sub>2</sub>R<sup>2a</sup>, OC(O)NR<sup>2</sup>R<sup>2a</sup>, C(O)NR<sup>2</sup>R<sup>2a</sup>,  
C(O)NR<sup>2</sup>(CH<sub>2</sub>)<sub>r</sub>OR<sup>2</sup>, SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>R<sup>2</sup>,

C(O)NR<sup>2</sup>SO<sub>2</sub>R<sup>2</sup>, SO<sub>2</sub>R<sup>2</sup>C(O)NR<sup>2</sup>, SO<sub>2</sub>NR<sup>2</sup>C(O)R<sup>2</sup>, C<sub>3</sub>-10 carbocycle substituted with 0-2 R<sup>4</sup>, and 4-10 membered heterocycle consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4</sup>, provided that R<sup>1b</sup> forms other than an O-O, N-halo, N-S, or N-CN bond;

R<sup>1c</sup> is selected from H, CH(CH<sub>2</sub>OR<sup>2</sup>)<sub>2</sub>, C(O)R<sup>2c</sup>, C(O)NR<sup>2</sup>R<sup>2a</sup>, S(O)R<sup>2</sup>, S(O)<sub>2</sub>R<sup>2</sup>, and SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>;

R<sup>1d</sup> is selected from C<sub>3</sub>-6 carbocycle substituted with 0-2 R<sup>4b</sup> and 5-10 membered heterocycle consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4b</sup>, provided that R<sup>1d</sup> forms other than an N-S bond;

R<sup>2</sup>, at each occurrence, is selected from H, CF<sub>3</sub>, C<sub>1</sub>-6 alkyl, benzyl, -(CH<sub>2</sub>)<sub>r</sub>-C<sub>3</sub>-10 carbocycle substituted with 0-2 R<sup>4b</sup>, and -(CH<sub>2</sub>)<sub>r</sub>-5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4b</sup>;

R<sup>2a</sup>, at each occurrence, is selected from H, CF<sub>3</sub>, C<sub>1</sub>-6 alkyl, benzyl, -(CH<sub>2</sub>)<sub>r</sub>-C<sub>3</sub>-10 carbocycle substituted with 0-2 R<sup>4b</sup>, and -(CH<sub>2</sub>)<sub>r</sub>-5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4b</sup>;

alternatively, R<sup>2</sup> and R<sup>2a</sup>, together with the atom to which they are attached, combine to form a 5-8 membered saturated, partially saturated or unsaturated ring substituted with 0-2 R<sup>4b</sup> and consisting of: 0-1 additional heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub>;

$R^{2b}$ , at each occurrence, is selected from  $CF_3$ ,  $C_{1-4}$  alkoxy substituted with 0-2  $R^{4b}$ ,  $C_{1-6}$  alkyl substituted with 0-2  $R^{4b}$ ,  $-(CH_2)_r-C_{3-10}$  carbocycle substituted with 0-2  $R^{4b}$ , and  $-(CH_2)_r-5-10$  membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and  $S(O)_p$  and substituted with 0-2  $R^{4b}$ ;

$R^{2c}$ , at each occurrence, is selected from  $CF_3$ , OH,  $C_{1-4}$  alkoxy,  $C_{1-6}$  alkyl,  $-(CH_2)_r-C_{3-10}$  carbocycle substituted with 0-2  $R^{4b}$ , and  $-(CH_2)_r-5-10$  membered heterocycle containing from 1-4 heteroatoms selected from the group consisting of N, O, and  $S(O)_p$  and substituted with 0-2  $R^{4b}$ ;

$R^3$ , at each occurrence, is selected from H,  $CH_3$ ,  $CH_2CH_3$ ,  $CH_2CH_2CH_3$ ,  $CH(CH_3)_2$ ,  $CH_2CH_2CH_2CH_3$ ,  $CH_2CH(CH_3)_2$ ,  $CH(CH_3)CH_2CH_3$ ,  $C(CH_3)_3$ , benzyl, and phenyl;

$R^{3a}$ , at each occurrence, is selected from H,  $CH_3$ ,  $CH_2CH_3$ ,  $CH_2CH_2CH_3$ ,  $CH(CH_3)_2$ ,  $CH_2CH_2CH_2CH_3$ ,  $CH_2CH(CH_3)_2$ ,  $CH(CH_3)CH_2CH_3$ ,  $C(CH_3)_3$ , benzyl, and phenyl;

alternatively,  $R^3$  and  $R^{3a}$ , together with the nitrogen atom to which they are attached, combine to form a 5 or 6 membered saturated, partially unsaturated, or unsaturated ring consisting of: carbon atoms, the nitrogen atom to which  $R^3$  and  $R^{3a}$  are attached, and 0-1 additional heteroatoms selected from the group consisting of N, O, and  $S(O)_p$ ;

$R^{3c}$ , at each occurrence, is selected from  $CH_3$ ,  $CH_2CH_3$ ,  $CH_2CH_2CH_3$ ,  $CH(CH_3)_2$ ,  $CH_2CH_2CH_2CH_3$ ,  $CH_2CH(CH_3)_2$ ,  $CH(CH_3)CH_2CH_3$ ,  $C(CH_3)_3$ , benzyl, and phenyl;

$R^{3d}$ , at each occurrence, is selected from H,  $CH_3$ ,  $CH_2CH_3$ ,  $CH_2CH_2CH_3$ ,  $CH(CH_3)_2$ ,  $CH_2CH_2CH_2CH_3$ ,  $CH_2CH(CH_3)_2$ ,  $CH(CH_3)CH_2CH_3$ ,  $C_{1-4}$  alkyl-phenyl, and  $C(=O)R^{3c}$ ;

$R^4$ , at each occurrence, is selected from H, =O,  $(CR^{3a})_rOR^2$ , F, Cl, Br, I,  $C_{1-4}$  alkyl,  $(CR^{3a})_rCN$ ,  $(CR^{3a})_rNO_2$ ,  $(CR^{3a})_rNR^2R^2a$ ,  $(CR^{3a})_rC(O)R^2c$ ,  $(CR^{3a})_rNR^2C(O)R^2b$ ,  $(CR^{3a})_rC(O)NR^2R^2a$ ,  $(CR^{3a})_rNR^3(CR^{3a})_rC(O)NR^3R^3a$ ,  $(CR^{3a})_rNR^3(CR^{3a})_rC(O)OR^3$ ,  $(CR^{3a})_rNR^3(CR^{3a})_rNR^3R^3a$ ,  $(CR^{3a})_rNR^3(CR^{3a})_rNR^3C(O)R^3a$ ,  $(CR^{3a})_rNR^3(CR^{3a})_rNR^3SO_2R^3a$ ,  $(CR^{3a})_rNR^2C(O)NR^2R^2a$ ,  $(CR^{3a})_rC(=NR^2)NR^2R^2a$ ,  $(CR^{3a})_rC(=NS(O)_2R^5)NR^2R^2a$ ,  $(CR^{3a})_rNHC(=NR^2)NR^2R^2a$ ,  $(CR^{3a})_rC(O)NHC(=NR^2)NR^2R^2a$ ,  $(CR^{3a})_rSO_2NR^2R^2a$ ,  $(CR^{3a})_rNR^2SO_2NR^2R^2a$ ,  $(CR^{3a})_rNR^2SO_2-C_{1-4}$  alkyl,  $(CR^{3a})_rNR^2SO_2R^5$ ,  $(CR^{3a})_rS(O)_pR^5a$ ,  $(CR^{3a})_r(CF_2)_rCF_3$ ,  $NHCH_2R^{1c}$ ,  $OCH_2R^{1c}$ ,  $SCH_2R^{1c}$ ,  $NH(CH_2)_2(CH_2)_tR^{1b}$ ,  $O(CH_2)_2(CH_2)_tR^{1b}$ ,  $S(CH_2)_2(CH_2)_tR^{1b}$ ,  $(CR^{3a})_r-3-10$  membered carbocycle substituted with 0-1  $R^5$ , and a  $(CR^{3a})_r-5-10$  membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and  $S(O)_p$  and substituted with 0-1  $R^5$ ;

$R^{4a}$ , at each occurrence, is selected from H, =O,  $(CR^{3a})_rOR^2$ ,  $(CR^{3a})_rF$ ,  $(CR^{3a})_rBr$ ,  $(CR^{3a})_rCl$ ,  $C_{1-4}$  alkyl,  $(CR^{3a})_rCN$ ,  $(CR^{3a})_rNO_2$ ,

(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>NR<sup>2</sup>R<sup>2a</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>C(O)R<sup>2c</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>NR<sup>2</sup>C(O)R<sup>2b</sup>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>C(O)NR<sup>2</sup>R<sup>2a</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>N=CHOR<sup>3</sup>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>C(O)NH(CH<sub>2</sub>)<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>NR<sup>2</sup>C(O)NR<sup>2</sup>R<sup>2a</sup>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>C(=NR<sup>2</sup>)NR<sup>2</sup>R<sup>2a</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>NHC(=NR<sup>2</sup>)NR<sup>2</sup>R<sup>2a</sup>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>NR<sup>2</sup>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>NR<sup>2</sup>SO<sub>2</sub>-C<sub>1-4</sub>  
alkyl, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>C(O)NHSO<sub>2</sub>-C<sub>1-4</sub> alkyl, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>NR<sup>2</sup>SO<sub>2</sub>R<sup>5</sup>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>S(O)<sub>p</sub>R<sup>5a</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>(CF<sub>2</sub>)<sub>r</sub>CF<sub>3</sub>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>-5-6 membered  
carbocycle substituted with 0-1 R<sup>5</sup>, and a (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>-5-6 membered heterocycle  
consisting of: carbon atoms and 1-4 heteroatoms selected from the group  
consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-1 R<sup>5</sup>;

R<sup>4b</sup>, at each occurrence, is selected from H, =O, (CH<sub>2</sub>)<sub>r</sub>OR<sup>3</sup>, (CH<sub>2</sub>)<sub>r</sub>F, (CH<sub>2</sub>)<sub>r</sub>Cl,  
(CH<sub>2</sub>)<sub>r</sub>Br, (CH<sub>2</sub>)<sub>r</sub>I, C<sub>1-4</sub> alkyl, (CH<sub>2</sub>)<sub>r</sub>CN, (CH<sub>2</sub>)<sub>r</sub>NO<sub>2</sub>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>R<sup>3a</sup>,  
(CH<sub>2</sub>)<sub>r</sub>C(O)R<sup>3</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)OR<sup>3c</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>C(O)R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>-C(O)NR<sup>3</sup>R<sup>3a</sup>,  
(CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>C(O)NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>C(=NR<sup>3</sup>)NR<sup>3</sup>R<sup>3a</sup>,  
(CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>C(=NR<sup>3</sup>)NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>SO<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>SO<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>,  
(CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>SO<sub>2</sub>CF<sub>3</sub>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>SO<sub>2</sub>-phenyl,  
(CH<sub>2</sub>)<sub>r</sub>S(O)<sub>p</sub>CF<sub>3</sub>, (CH<sub>2</sub>)<sub>r</sub>S(O)<sub>p</sub>-C<sub>1-4</sub> alkyl, (CH<sub>2</sub>)<sub>r</sub>S(O)<sub>p</sub>-phenyl,  
(CH<sub>2</sub>)<sub>r</sub>(CF<sub>2</sub>)<sub>r</sub>CF<sub>3</sub>, (CH<sub>2</sub>)<sub>r</sub>-3-10 membered carbocycle substituted with 0-1 R<sup>3</sup>,  
and a (CH<sub>2</sub>)<sub>r</sub>-5-10 membered heterocycle consisting of: carbon atoms and 1-4  
heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and  
substituted with 0-1 R<sup>3</sup>;

R<sup>4c</sup>, at each occurrence, is selected from H, C<sub>1-4</sub> alkyl (CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>OR<sup>2</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>F,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>Br, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>Cl, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>CN, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>NO<sub>2</sub>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>NR<sup>2</sup>R<sup>2a</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>C(O)R<sup>2c</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>NR<sup>2</sup>C(O)R<sup>2b</sup>,

(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>C(O)NR<sup>2</sup>R<sup>2a</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>N=CHOR<sup>3</sup>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>C(O)NH(CH<sub>2</sub>)<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>NR<sup>2</sup>C(O)NR<sup>2</sup>R<sup>2a</sup>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>C(=NR<sup>2</sup>)NR<sup>2</sup>R<sup>2a</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>NHC(=NR<sup>2</sup>)NR<sup>2</sup>R<sup>2a</sup>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>NR<sup>2</sup>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>NR<sup>2</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>C(O)NHSO<sub>2</sub>-C<sub>1-4</sub> alkyl,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r1</sub>NR<sup>2</sup>SO<sub>2</sub>R<sup>5</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>S(O)<sub>p</sub>R<sup>5a</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>(CF<sub>2</sub>)<sub>r</sub>CF<sub>3</sub>,  
(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>-5-6 membered carbocycle substituted with 0-1 R<sup>5</sup>, and a (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>-  
5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms  
selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-1  
R<sup>5</sup>;

R<sup>5</sup>, at each occurrence, is selected from H, C<sub>1-6</sub> alkyl, =O, (CH<sub>2</sub>)<sub>r</sub>OR<sup>3</sup>, F, Cl, Br, I, -  
CN, NO<sub>2</sub>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)R<sup>3</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)OR<sup>3c</sup>,  
(CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>C(O)R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>C(O)NR<sup>3</sup>R<sup>3a</sup>,  
(CH<sub>2</sub>)<sub>r</sub>CH(=NOR<sup>3d</sup>), (CH<sub>2</sub>)<sub>r</sub>C(=NR<sup>3</sup>)NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>C(=NR<sup>3</sup>)NR<sup>3</sup>R<sup>3a</sup>,  
(CH<sub>2</sub>)<sub>r</sub>SO<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>SO<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl,  
(CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>SO<sub>2</sub>CF<sub>3</sub>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>SO<sub>2</sub>-phenyl, (CH<sub>2</sub>)<sub>r</sub>S(O)<sub>p</sub>CF<sub>3</sub>,  
(CH<sub>2</sub>)<sub>r</sub>S(O)<sub>p</sub>-C<sub>1-4</sub> alkyl, (CH<sub>2</sub>)<sub>r</sub>S(O)<sub>p</sub>-phenyl, (CF<sub>2</sub>)<sub>r</sub>CF<sub>3</sub>, phenyl substituted  
with 0-2 R<sup>6</sup>, naphthyl substituted with 0-2 R<sup>6</sup>, and benzyl substituted with 0-2  
R<sup>6</sup>;

R<sup>5a</sup>, at each occurrence, is selected from C<sub>1-6</sub> alkyl, (CH<sub>2</sub>)<sub>r</sub>OR<sup>3</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>R<sup>3a</sup>,  
(CH<sub>2</sub>)<sub>r</sub>C(O)R<sup>3</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)OR<sup>3c</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>C(O)R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)NR<sup>3</sup>R<sup>3a</sup>,  
(CF<sub>2</sub>)<sub>r</sub>CF<sub>3</sub>, phenyl substituted with 0-2 R<sup>6</sup>, naphthyl substituted with 0-2 R<sup>6</sup>, and  
benzyl substituted with 0-2 R<sup>6</sup>, provided that R<sup>5a</sup> does not form a S-N or S(O)<sub>p</sub>-  
C(O) bond;

R<sup>6</sup>, at each occurrence, is selected from H, OH, (CH<sub>2</sub>)<sub>r</sub>OR<sup>2</sup>, halo, C<sub>1-4</sub> alkyl, CN, NO<sub>2</sub>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>2</sup>R<sup>2a</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)R<sup>2b</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>, NR<sup>2</sup>C(O)NR<sup>2</sup>R<sup>2a</sup>, C(=NH)NH<sub>2</sub>, NHC(=NH)NH<sub>2</sub>, SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, and NR<sup>2</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl;

R<sup>7</sup>, at each occurrence, is selected from H, OH, C<sub>1-6</sub> alkyl, C<sub>1-6</sub> alkyl-C(O)-, C<sub>1-6</sub> alkyl-O-, (CH<sub>2</sub>)<sub>n</sub>-phenyl, C<sub>1-6</sub> alkyl-OC(O)-, C<sub>6-10</sub> aryl-O-, C<sub>6-10</sub> aryl-OC(O)-, C<sub>6-10</sub> aryl-CH<sub>2</sub>-C(O)-, C<sub>1-4</sub> alkyl-C(O)O-C<sub>1-4</sub> alkyl-OC(O)-, C<sub>6-10</sub> aryl-C(O)O-C<sub>1-4</sub> alkyl-OC(O)-, C<sub>1-6</sub> alkyl-NH<sub>2</sub>-C(O)-, phenyl-NH<sub>2</sub>-C(O)-, and phenyl C<sub>0-4</sub> alkyl-C(O)-;

R<sup>8</sup>, at each occurrence, is selected from H, C<sub>1-6</sub> alkyl, and (CH<sub>2</sub>)<sub>n</sub>-phenyl;

alternatively, R<sup>7</sup> and R<sup>8</sup>, when attached to the same nitrogen, combine to form a 5-10 membered heterocyclic ring consisting of carbon atoms and 0-2 additional heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub>;

R<sup>9</sup>, at each occurrence, is selected from H, C<sub>1-6</sub> alkyl, and (CH<sub>2</sub>)<sub>n</sub>-phenyl;

n, at each occurrence, is selected from 0, 1, 2, and 3;

p, at each occurrence, is selected from 0, 1, and 2;

r, at each occurrence, is selected from 0, 1, 2, 3, 4, 5, and 6;

r1, at each occurrence, is selected from 1, 2, 3, 4, 5, and 6;

t, at each occurrence, is selected from 0, 1, 2, and 3.

2. (Currently Amended) A compound according to Claim 1, wherein:

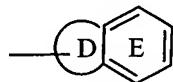
~~one of P and M<sub>1</sub> is G and the other A-B;~~

~~M is 3-8 membered linear chain consisting of: carbon atoms, 1-3 carbonyl groups, 0-1 thiocarbonyl groups, and 1-3 heteroatoms selected from O, S(O)<sub>p</sub>, and N; and M is substituted with 0-3 R<sup>1a</sup> and 0-2 R<sup>2</sup> and there are 0-1 double bonds, provided that other than an S-S, S-O, or O-O bond is present in M;~~

G is a group of formula IIa or IIb:



IIa



IIb

ring D, including the two atoms of Ring E to which it is attached, is a 5-6 membered ring consisting of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub>;

ring D is substituted with 0-2 R and there are 0-3 ring double bonds;

E is selected from phenyl, pyridyl, pyrimidyl, pyrazinyl, and pyridazinyl, and is substituted with 1-2 R;

alternatively, ring D is absent, and ring E is selected from phenyl, pyridyl, pyridazinyl, pyrimidyl, and thienyl, and ring E is substituted with 1-2 R;

alternatively, ring D is absent, ring E is selected from phenyl, pyridyl, and thienyl, and ring E is substituted with 1 R and with a 5 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O,

and S(O)p, wherein the 5 membered heterocycle is substituted with 0-1 carbonyls and 1-2 R and there are 0-3 ring double bonds;

R is selected from H, C<sub>1-4</sub> alkyl, F, Cl, OH, OCH<sub>3</sub>, OCH<sub>2</sub>CH<sub>3</sub>, OCH(CH<sub>3</sub>)<sub>2</sub>, CN, C(=NH)NH<sub>2</sub>, C(=NH)NHOH, C(=NH)NHOCH<sub>3</sub>, NH<sub>2</sub>, NH(C<sub>1-3</sub> alkyl), N(C<sub>1-3</sub> alkyl)<sub>2</sub>, C(=NH)NH<sub>2</sub>, CH<sub>2</sub>NH<sub>2</sub>, CH<sub>2</sub>NH(C<sub>1-3</sub> alkyl), CH<sub>2</sub>N(C<sub>1-3</sub> alkyl)<sub>2</sub>, (CR<sup>8</sup>R<sup>9</sup>)<sub>t</sub>NR<sup>7</sup>R<sup>8</sup>, C(O)NR<sup>7</sup>R<sup>8</sup>, CH<sub>2</sub>C(O)NR<sup>7</sup>R<sup>8</sup>, S(O)<sub>2</sub>R<sup>3</sup>, S(O)pNR<sup>7</sup>R<sup>8</sup>, CH<sub>2</sub>S(O)pNR<sup>7</sup>R<sup>8</sup>, and OCF<sub>3</sub>;

alternatively, when 2 R groups are attached to adjacent atoms, they combine to form methylenedioxy or ethylenedioxy;

A is selected from:

C<sub>5-10</sub> carbocycle substituted with 0-2 R<sup>4</sup>, and  
5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)p and substituted with 0-2 R<sup>4</sup>;

B is ; provided that Z and B are attached to different atoms on A and that the A-X-N moiety forms other than a N-N-N group;

ring Q is a 4-7 membered monocyclic or tricyclic ring consisting of, in addition to the N-Q<sub>1</sub> group shown, carbon atoms and 0-2 heteroatoms selected from NR<sup>4c</sup>, O, S, S(O), and S(O)<sub>2</sub>, wherein:

0-2 double bonds are present within the ring and the ring is substituted with 0-2 R<sup>4a</sup>;

alternatively, ring Q is a 4-7 membered ring to which another ring is fused, wherein:

the 4-7 membered ring consists of, in addition to the shown amide group, carbon atoms and 0-2 heteroatoms selected from NR<sup>4c</sup>, O, S, S(O), and S(O)<sub>2</sub> and 0-1 double bonds are present within the ring;

the fusion ring is phenyl or a 5-6 membered heteroaromatic consisting of carbon atoms and 1-2 heteroatoms selected from NR<sup>4c</sup>, O, and S; ring Q, which includes the 4-7 membered ring and the fusion ring, is substituted with 0-3 R<sup>4a</sup>;

X is absent or is selected from -(CR<sup>2</sup>R<sup>2a</sup>)<sub>1-4</sub>-, -C(O)-, -C(O)CR<sup>2</sup>R<sup>2a</sup>-, -CR<sup>2</sup>R<sup>2a</sup>C(O)-, -S(O)<sub>2</sub>-, -S(O)<sub>2</sub>CR<sup>2</sup>R<sup>2a</sup>-, -CR<sup>2</sup>R<sup>2a</sup>S(O)<sub>2</sub>-, -NR<sup>2</sup>S(O)<sub>2</sub>-, -NR<sup>2</sup>CR<sup>2</sup>R<sup>2a</sup>-, and -OCR<sup>2</sup>R<sup>2a</sup>-;

R<sup>1a</sup>, at each occurrence, is selected from H, -(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>-R<sup>1b</sup>, -(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>-O-(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>-R<sup>1b</sup>, -C<sub>2-6</sub> alkenylene-R<sup>1b</sup>, -C<sub>2-6</sub> alkynylene-R<sup>1b</sup>, -(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>-C(=NR<sup>1b</sup>)NR<sup>3</sup>R<sup>1b</sup>, NR<sup>3</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1c</sup>, O(CR<sup>3</sup>R<sup>3a</sup>)<sub>t</sub>R<sup>1c</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>SCR<sup>3</sup>R<sup>3a</sup>R<sup>1c</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>NR<sup>3</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1b</sup>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>C(O)NR<sup>2</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1b</sup>, CO<sub>2</sub>(CR<sup>3</sup>R<sup>3a</sup>)<sub>t</sub>R<sup>1b</sup>, O(CR<sup>3</sup>R<sup>3a</sup>)<sub>t</sub>R<sup>1b</sup>, S(O)<sub>p</sub>(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>, O(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>, NR<sup>3</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>, OC(O)NR<sup>3</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>, NR<sup>3</sup>C(O)NR<sup>3</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>, NR<sup>3</sup>C(O)O(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>, and NR<sup>3</sup>C(O)(CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>R<sup>1d</sup>, provided that R<sup>1a</sup> forms other than an N-halo, N-S, O-O, or N-CN bond;

alternatively, when two R<sup>1a</sup> groups are attached to the same carbon atom, together with the carbon atom to which they are attached they form a 3-10 membered carbocyclic or heterocyclic ring consisting of: carbon atoms and 0-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub>, this ring being substituted with 0-2 R<sup>4</sup> and 0-3 ring double bonds;

R<sup>1b</sup> is selected from H, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, F, Cl, Br, I, -CN, -CHO, CF<sub>3</sub>, (CR<sup>3</sup>R<sup>3a</sup>)<sub>r</sub>OR<sup>2</sup>, NR<sup>2</sup>R<sup>2a</sup>, C(O)R<sup>2b</sup>, CO<sub>2</sub>R<sup>2b</sup>, OC(O)R<sup>2</sup>, CO<sub>2</sub>R<sup>2a</sup>, S(O)<sub>p</sub>R<sup>2</sup>, NR<sup>2</sup>(CH<sub>2</sub>)<sub>r</sub>OR<sup>2</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>, NR<sup>2</sup>C(O)NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>C(O)<sub>2</sub>R<sup>2a</sup>, OC(O)NR<sup>2</sup>R<sup>2a</sup>, C(O)NR<sup>2</sup>R<sup>2a</sup>, C(O)NR<sup>2</sup>(CH<sub>2</sub>)<sub>r</sub>OR<sup>2</sup>, SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>R<sup>2</sup>, C(O)NR<sup>2</sup>SO<sub>2</sub>R<sup>2</sup>, SO<sub>2</sub>NR<sup>2</sup>C(O)R<sup>2</sup>, C<sub>3-10</sub> carbocycle substituted with 0-2 R<sup>4</sup>, and 4-10 membered heterocycle consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4</sup>, provided that R<sup>1b</sup> forms other than an O-O, N-halo, N-S, or N-CN bond;

R<sup>1c</sup> is selected from H, CH(CH<sub>2</sub>OR<sup>2</sup>)<sub>2</sub>, C(O)R<sup>2c</sup>, C(O)NR<sup>2</sup>R<sup>2a</sup>, S(O)R<sup>2</sup>, S(O)<sub>2</sub>R<sup>2</sup>, and SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>;

R<sup>2</sup>, at each occurrence, is selected from H, CF<sub>3</sub>, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH(CH<sub>3</sub>)<sub>2</sub>, CH(CH<sub>3</sub>)CH<sub>2</sub>CH<sub>3</sub>, C(CH<sub>3</sub>)<sub>3</sub>, benzyl, C<sub>5-6</sub> carbocycle substituted with 0-2 R<sup>4b</sup>, a C<sub>5-6</sub> carbocyclic-CH<sub>2</sub>-group substituted with 0-2 R<sup>4b</sup>, and 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4b</sup>;

R<sup>2a</sup>, at each occurrence, is selected from H, CF<sub>3</sub>, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH(CH<sub>3</sub>)<sub>2</sub>, CH(CH<sub>3</sub>)CH<sub>2</sub>CH<sub>3</sub>, C(CH<sub>3</sub>)<sub>3</sub>, benzyl, C<sub>5-6</sub> carbocycle substituted with 0-2 R<sup>4b</sup>, and 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4b</sup>;

alternatively,  $R^2$  and  $R^{2a}$ , together with the atom to which they are attached, combine to form a 5 or 6 membered saturated, partially saturated or unsaturated ring substituted with 0-2  $R^{4b}$  and consisting of: 0-1 additional heteroatoms selected from the group consisting of N, O, and S(O)p;

$R^{2b}$ , at each occurrence, is selected from  $CF_3$ , C<sub>1-4</sub> alkoxy,  $CH_3$ ,  $CH_2CH_3$ ,  $CH_2CH_2CH_3$ ,  $CH(CH_3)_2$ ,  $CH_2CH_2CH_2CH_3$ ,  $CH_2CH(CH_3)_2$ ,  $CH(CH_3)CH_2CH_3$ ,  $C(CH_3)_3$ , benzyl, C<sub>5-6</sub> carbocycle substituted with 0-2  $R^{4b}$ , and 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)p and substituted with 0-2  $R^{4b}$ ;

$R^{2c}$ , at each occurrence, is selected from  $CF_3$ , OH, C<sub>1-4</sub> alkoxy,  $CH_3$ ,  $CH_2CH_3$ ,  $CH_2CH_2CH_3$ ,  $CH(CH_3)_2$ ,  $CH_2CH_2CH_2CH_3$ ,  $CH_2CH(CH_3)_2$ ,  $CH(CH_3)CH_2CH_3$ ,  $C(CH_3)_3$ , benzyl, C<sub>5-6</sub> carbocycle substituted with 0-2  $R^{4b}$ , and 5-6 membered heterocycle containing from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)p and substituted with 0-2  $R^{4b}$ ;

$R^3$ , at each occurrence, is selected from H,  $CH_3$ ,  $CH_2CH_3$ ,  $CH_2CH_2CH_3$ ,  $CH(CH_3)_2$ , benzyl, and phenyl;

$R^{3a}$ , at each occurrence, is selected from H,  $CH_3$ ,  $CH_2CH_3$ ,  $CH_2CH_2CH_3$ ,  $CH(CH_3)_2$ , benzyl, and phenyl;

alternatively,  $R^3$  and  $R^{3a}$ , together with the nitrogen atom to which they are attached, combine to form a 5 or 6 membered saturated, partially unsaturated, or unsaturated ring consisting of: carbon atoms and the nitrogen atom to which  $R^3$  and  $R^{3a}$  are attached;

R<sup>3c</sup>, at each occurrence, is selected from CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, benzyl, and phenyl;

R<sup>3d</sup>, at each occurrence, is selected from H, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>2</sub>-phenyl, CH<sub>2</sub>CH<sub>2</sub>-phenyl, and C(=O)R<sup>3c</sup>;

R<sup>4</sup>, at each occurrence, is selected from H, =O, (CH<sub>2</sub>)<sub>r</sub>OR<sup>2</sup>, F, Cl, Br, I, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH(CH<sub>3</sub>)<sub>2</sub>, CH(CH<sub>3</sub>)CH<sub>2</sub>CH<sub>3</sub>, C(CH<sub>3</sub>)<sub>3</sub>, -CN, NO<sub>2</sub>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>2</sup>R<sup>2a</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)R<sup>2c</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>2</sup>C(O)R<sup>2b</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)NR<sup>2</sup>R<sup>2a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>(CH<sub>2</sub>)<sub>1-4</sub>C(O)NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>(CH<sub>2</sub>)<sub>1-4</sub>C(O)OR<sup>3</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>(CH<sub>2</sub>)<sub>1-4</sub>NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>(CH<sub>2</sub>)<sub>1-4</sub>NR<sup>3</sup>C(O)R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>(CH<sub>2</sub>)<sub>1-4</sub>NR<sup>3</sup>SO<sub>2</sub>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>2</sup>C(O)NR<sup>2</sup>R<sup>2a</sup>, (CH<sub>2</sub>)<sub>r</sub>C(=NR<sup>2</sup>)NR<sup>2</sup>R<sup>2a</sup>, (CH<sub>2</sub>)<sub>r</sub>NHC(=NR<sup>2</sup>)NR<sup>2</sup>R<sup>2a</sup>, (CH<sub>2</sub>)<sub>r</sub>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>2</sup>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>2</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, (CH<sub>2</sub>)<sub>r</sub>NR<sup>2</sup>SO<sub>2</sub>R<sup>5</sup>, (CH<sub>2</sub>)<sub>r</sub>S(O)<sub>p</sub>R<sup>5a</sup>, (CH<sub>2</sub>)<sub>r</sub>CF<sub>3</sub>, (CH<sub>2</sub>)<sub>r</sub>-3-7 membered carbocycle substituted with 0-1 R<sup>5</sup>, and a (CH<sub>2</sub>)<sub>r</sub>-5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-1 R<sup>5</sup>;

R<sup>4a</sup>, at each occurrence, is selected from H, =O, CH<sub>2</sub>OR<sup>2</sup>, OR<sup>2</sup>, CH<sub>2</sub>F, F, CH<sub>2</sub>Br, Br, CH<sub>2</sub>Cl, Cl, C<sub>1-4</sub> alkyl, CH<sub>2</sub>-CN, -CN, CH<sub>2</sub>NO<sub>2</sub>, NO<sub>2</sub>, CH<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>-C(O)R<sup>2c</sup>, C(O)R<sup>2c</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>C(O)NR<sup>2</sup>R<sup>2a</sup>, (CH<sub>2</sub>)<sub>r</sub>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, NR<sup>2</sup>SO<sub>2</sub>R<sup>5</sup>, (CH<sub>2</sub>)<sub>r</sub>S(O)<sub>p</sub>R<sup>5a</sup>, CH<sub>2</sub>CF<sub>3</sub>, CF<sub>3</sub>, CH<sub>2</sub>-5-6 membered carbocycle substituted with 0-1 R<sup>5</sup>, 5-6 membered carbocycle substituted with 0-1 R<sup>5</sup>, and a

CH<sub>2</sub>-5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-1 R<sup>5</sup>, and 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-1 R<sup>5</sup>;

R<sup>4b</sup>, at each occurrence, is selected from H, =O, OR<sup>3</sup>, (CH<sub>2</sub>)<sub>r</sub>OR<sup>3</sup>, F, Cl, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH(CH<sub>3</sub>)<sub>2</sub>, CH(CH<sub>3</sub>)CH<sub>2</sub>CH<sub>3</sub>, C(CH<sub>3</sub>)<sub>3</sub>, -CN, NO<sub>2</sub>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)R<sup>3</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)OR<sup>3c</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>C(O)R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>C(O)NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>C(=NR<sup>3</sup>)NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>C(=NR<sup>3</sup>)NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>SO<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>SO<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>SO<sub>2</sub>CF<sub>3</sub>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>SO<sub>2</sub>-phenyl, (CH<sub>2</sub>)<sub>r</sub>S(O)<sub>p</sub>CF<sub>3</sub>, (CH<sub>2</sub>)<sub>r</sub>S(O)<sub>p</sub>-C<sub>1-4</sub> alkyl, (CH<sub>2</sub>)<sub>r</sub>S(O)<sub>p</sub>-phenyl, and (CH<sub>2</sub>)<sub>r</sub>CF<sub>3</sub>;

R<sup>4c</sup>, at each occurrence, is selected from H, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH(CH<sub>3</sub>)<sub>2</sub>, CH(CH<sub>3</sub>)CH<sub>2</sub>CH<sub>3</sub>, C(CH<sub>3</sub>)<sub>3</sub>, CH<sub>2</sub>OR<sup>2</sup>, CH<sub>2</sub>F, CH<sub>2</sub>Br, CH<sub>2</sub>Cl, CH<sub>2</sub>CN, CH<sub>2</sub>NO<sub>2</sub>, CH<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, C(O)R<sup>2c</sup>, CH<sub>2</sub>C(O)R<sup>2c</sup>, CH<sub>2</sub>NR<sup>2</sup>C(O)R<sup>2b</sup>, C(O)NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>C(O)NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>NR<sup>2</sup>C(O)NR<sup>2</sup>R<sup>2a</sup>, SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>NR<sup>2</sup>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>NR<sup>2</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, C(O)NHSO<sub>2</sub>-C<sub>1-4</sub> alkyl, CH<sub>2</sub>C(O)NHSO<sub>2</sub>-C<sub>1-4</sub> alkyl, CH<sub>2</sub>NR<sup>2</sup>SO<sub>2</sub>R<sup>5</sup>, S(O)<sub>p</sub>R<sup>5a</sup>, CH<sub>2</sub>S(O)<sub>p</sub>R<sup>5a</sup>, CF<sub>3</sub>, CH<sub>2</sub>CF<sub>3</sub>, 5-6 membered carbocycle substituted with 0-1 R<sup>5</sup>, CH<sub>2</sub>5-6 membered carbocycle substituted with 0-1 R<sup>5</sup>, 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and

substituted with 0-1 R<sup>5</sup>, and a CH<sub>2</sub>5-6 membered heterocycle consisting of:  
carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O,  
and S(O)<sub>p</sub> and substituted with 0-1 R<sup>5</sup>;

R<sup>5</sup>, at each occurrence, is selected from H, =O, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>,  
CH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH(CH<sub>3</sub>)<sub>2</sub>, CH(CH<sub>3</sub>)CH<sub>2</sub>CH<sub>3</sub>, C(CH<sub>3</sub>)<sub>3</sub>,  
OR<sup>3</sup>, CH<sub>2</sub>OR<sup>3</sup>, F, Cl, -CN, NO<sub>2</sub>, NR<sup>3</sup>R<sup>3a</sup>, CH<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, C(O)R<sup>3</sup>, CH<sub>2</sub>C(O)R<sup>3</sup>,  
C(O)OR<sup>3c</sup>, CH<sub>2</sub>C(O)OR<sup>3c</sup>, NR<sup>3</sup>C(O)R<sup>3a</sup>, C(O)NR<sup>3</sup>R<sup>3a</sup>, NR<sup>3</sup>C(O)NR<sup>3</sup>R<sup>3a</sup>,  
CH(=NOR<sup>3d</sup>), C(=NR<sup>3</sup>)NR<sup>3</sup>R<sup>3a</sup>, NR<sup>3</sup>C(=NR<sup>3</sup>)NR<sup>3</sup>R<sup>3a</sup>, SO<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>,  
NR<sup>3</sup>SO<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, NR<sup>3</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, NR<sup>3</sup>SO<sub>2</sub>CF<sub>3</sub>, NR<sup>3</sup>SO<sub>2</sub>-phenyl,  
S(O)<sub>p</sub>CF<sub>3</sub>, S(O)<sub>p</sub>-C<sub>1-4</sub> alkyl, S(O)<sub>p</sub>-phenyl, CF<sub>3</sub>, phenyl substituted with 0-2 R<sup>6</sup>,  
naphthyl substituted with 0-2 R<sup>6</sup>, and benzyl substituted with 0-2 R<sup>6</sup>;

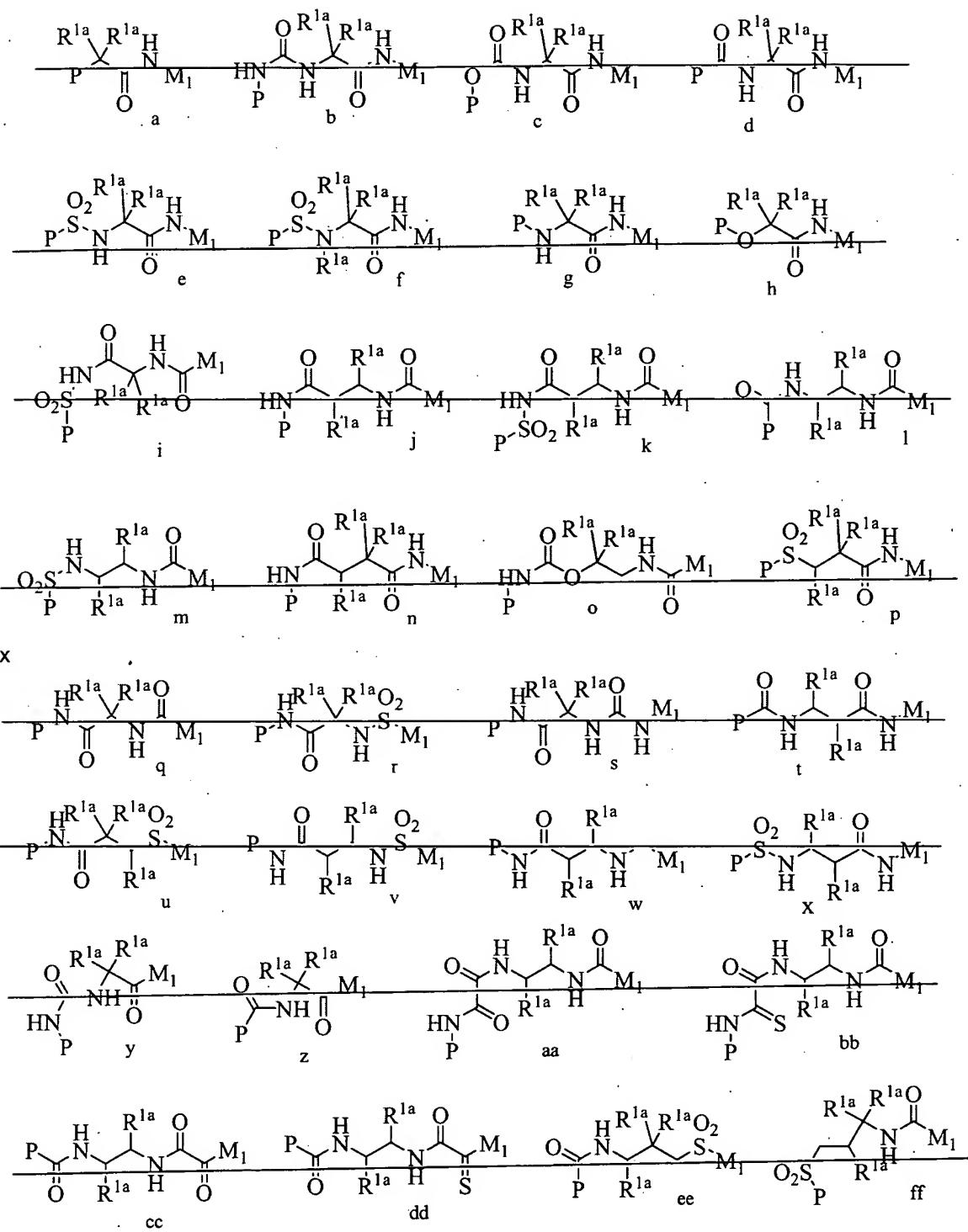
R<sup>6</sup>, at each occurrence, is selected from H, OH, OR<sup>2</sup>, F, Cl, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>,  
CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH(CH<sub>3</sub>)<sub>2</sub>,  
CH(CH<sub>3</sub>)CH<sub>2</sub>CH<sub>3</sub>, C(CH<sub>3</sub>)<sub>3</sub>, -CN, NO<sub>2</sub>, NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, C(O)R<sup>2b</sup>,  
CH<sub>2</sub>C(O)R<sup>2b</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>, NR<sup>2</sup>C(O)NR<sup>2</sup>R<sup>2a</sup>, C(=NH)NH<sub>2</sub>, NHC(=NH)NH<sub>2</sub>,  
SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, and NR<sup>2</sup>SO<sub>2</sub>C<sub>1-4</sub> alkyl;

r, at each occurrence, is selected from 0, 1, 2, and 3;

r<sub>1</sub>, at each occurrence, is selected from 1, 2, and 3; and,

t, at each occurrence, is selected from 0, 1, and 2.

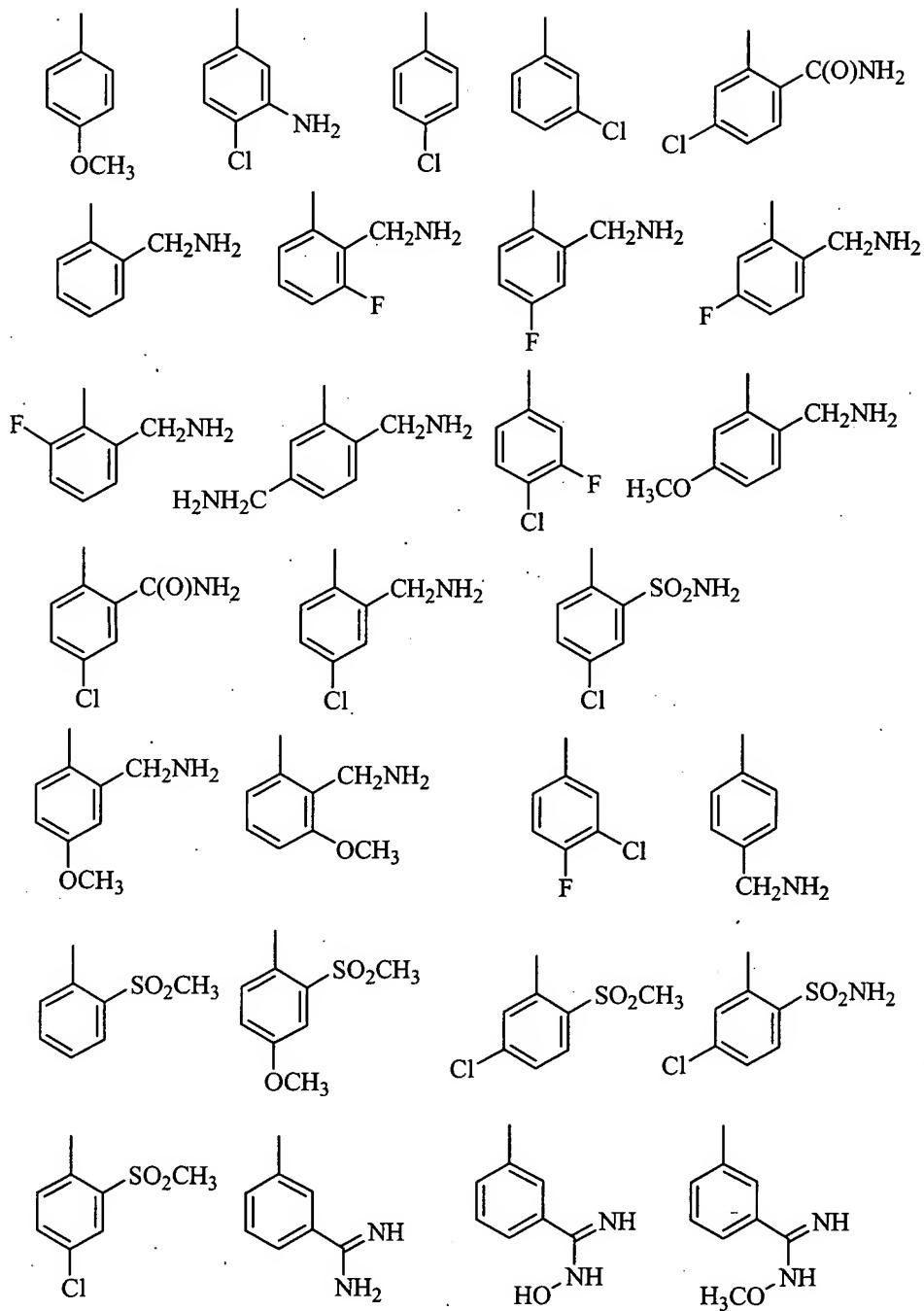
3. (Currently Amended) A compound according to Claim 2, wherein ~~the compound is selected from compounds a-ff~~:

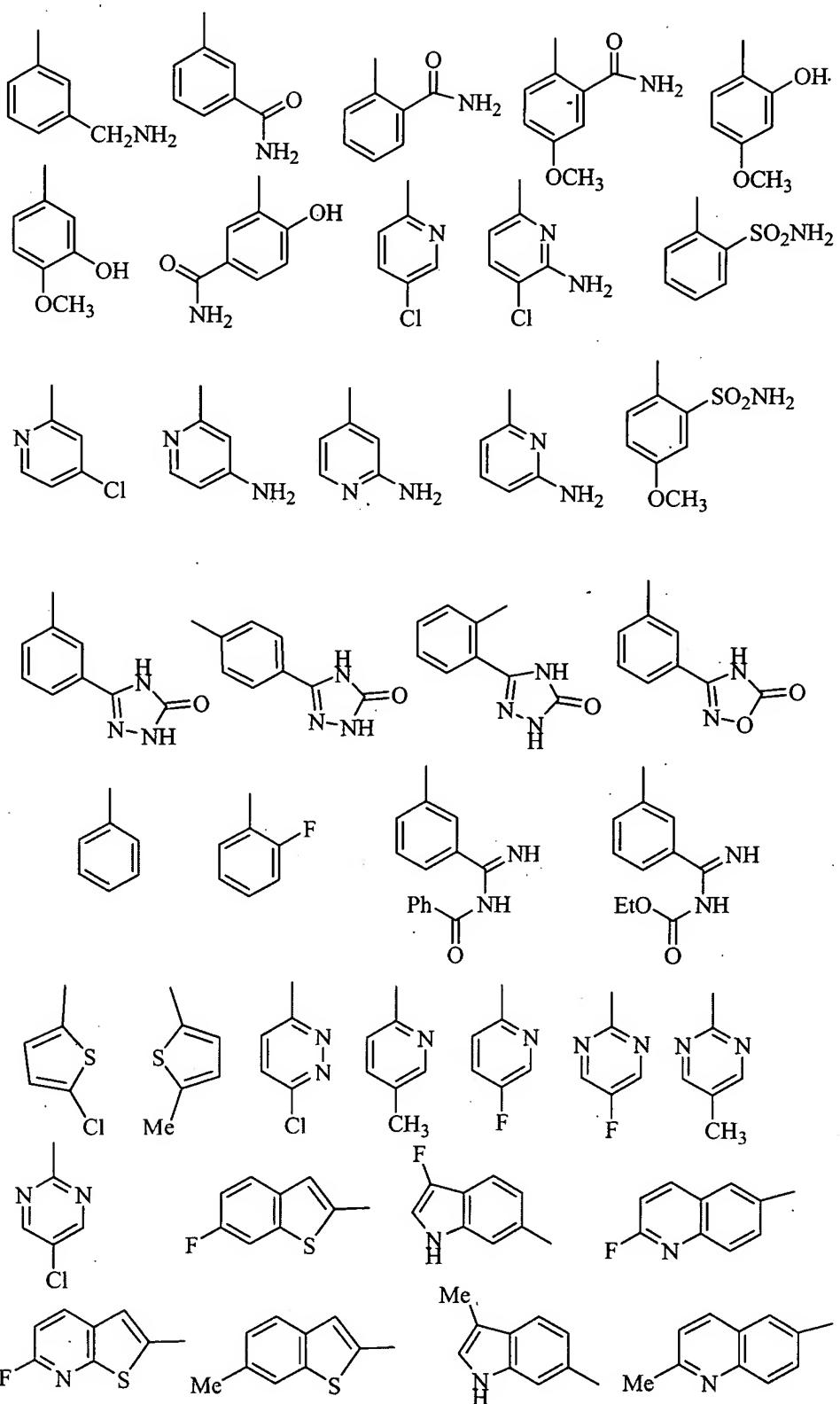


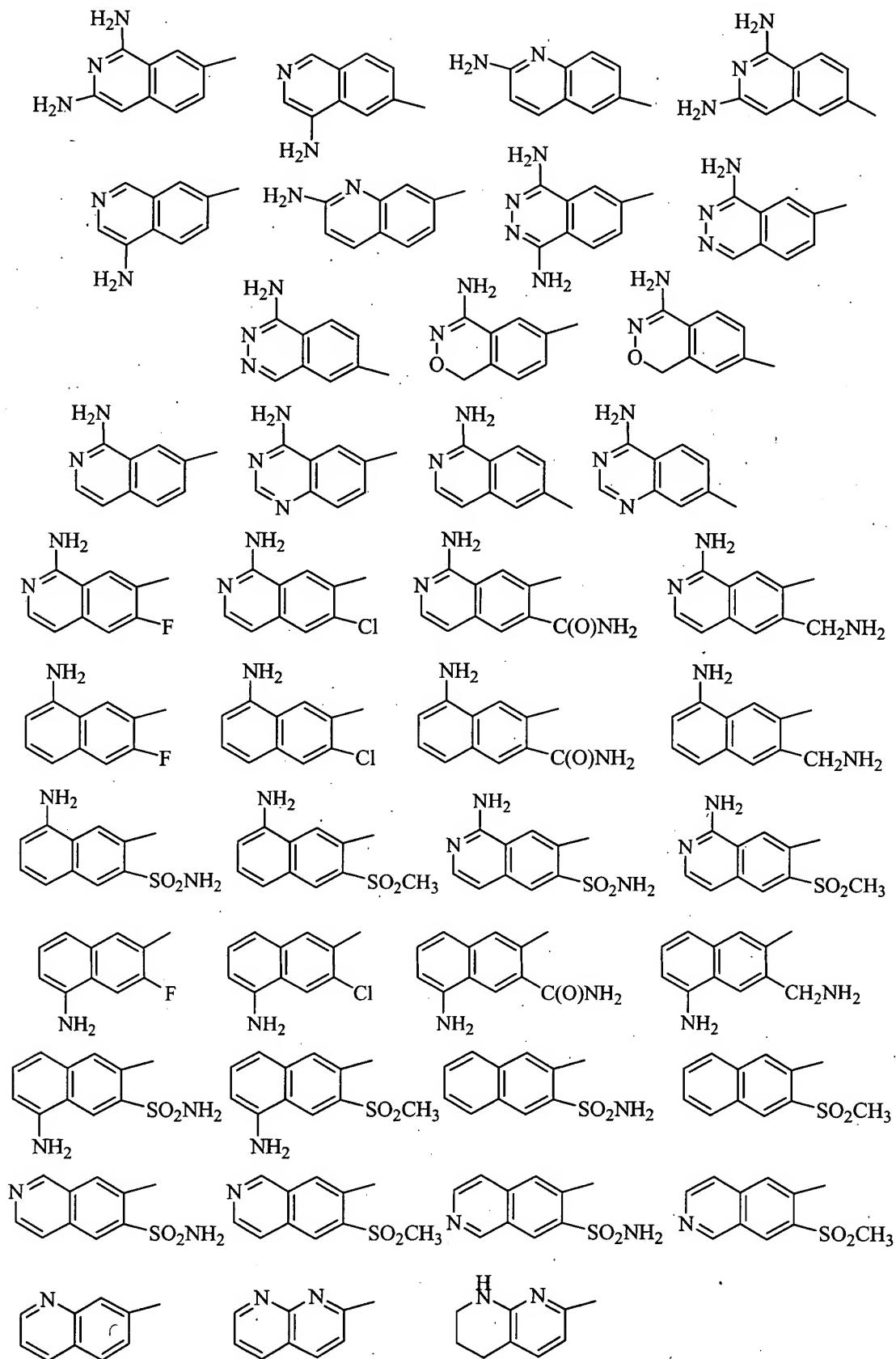
wherein:

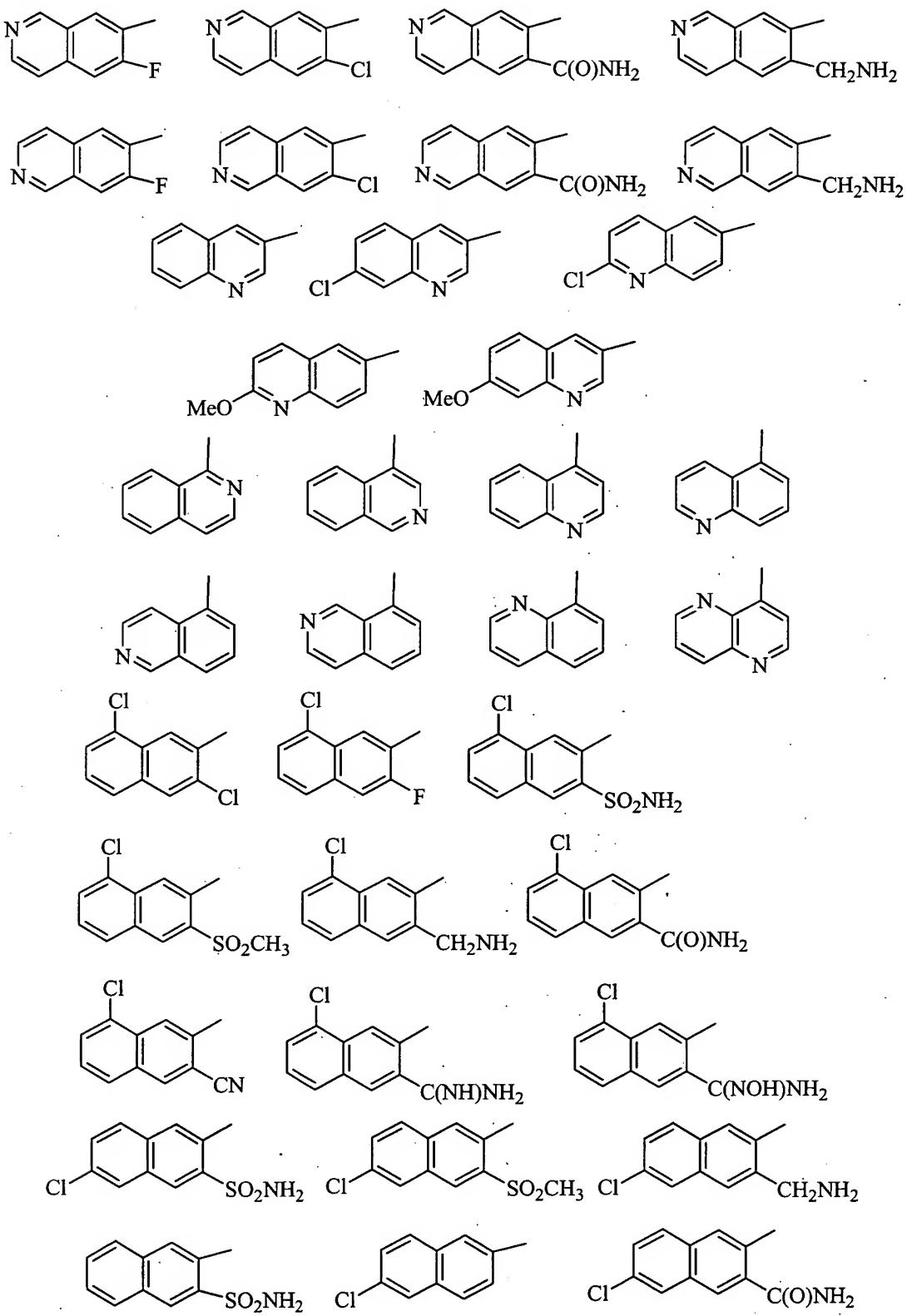
**one of P and M<sub>1</sub> is G and the other A-B;**

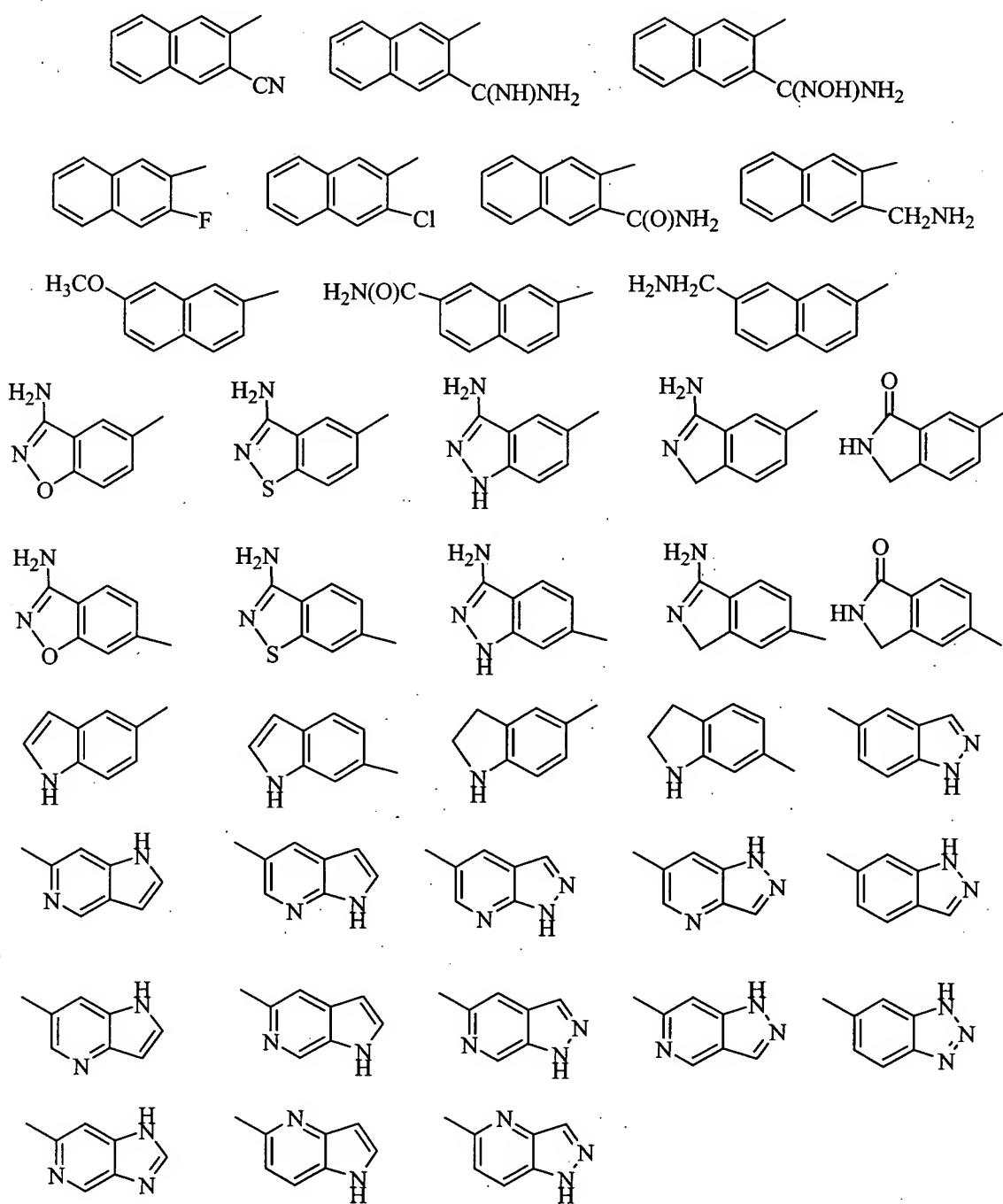
G is selected from the group:

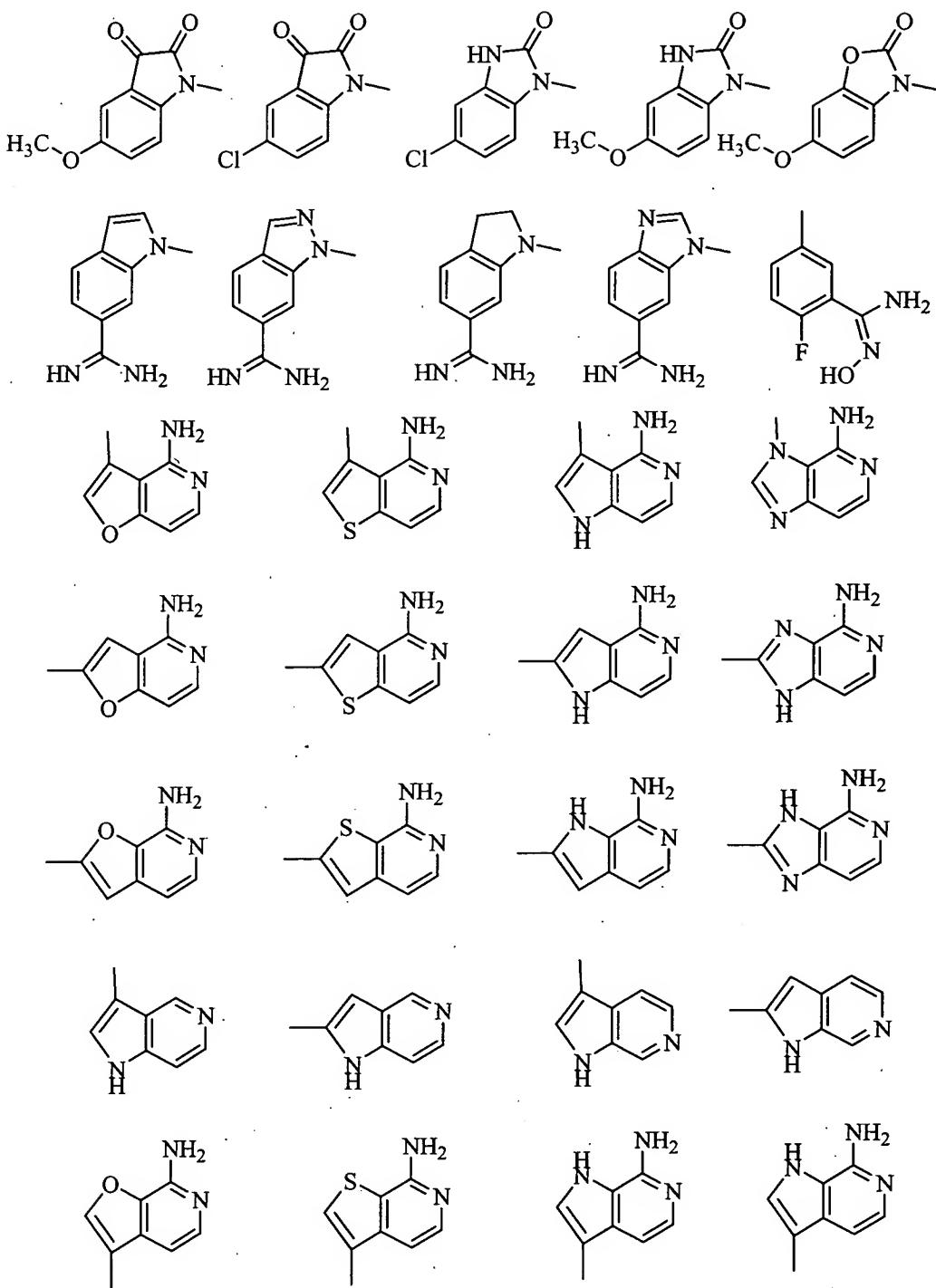


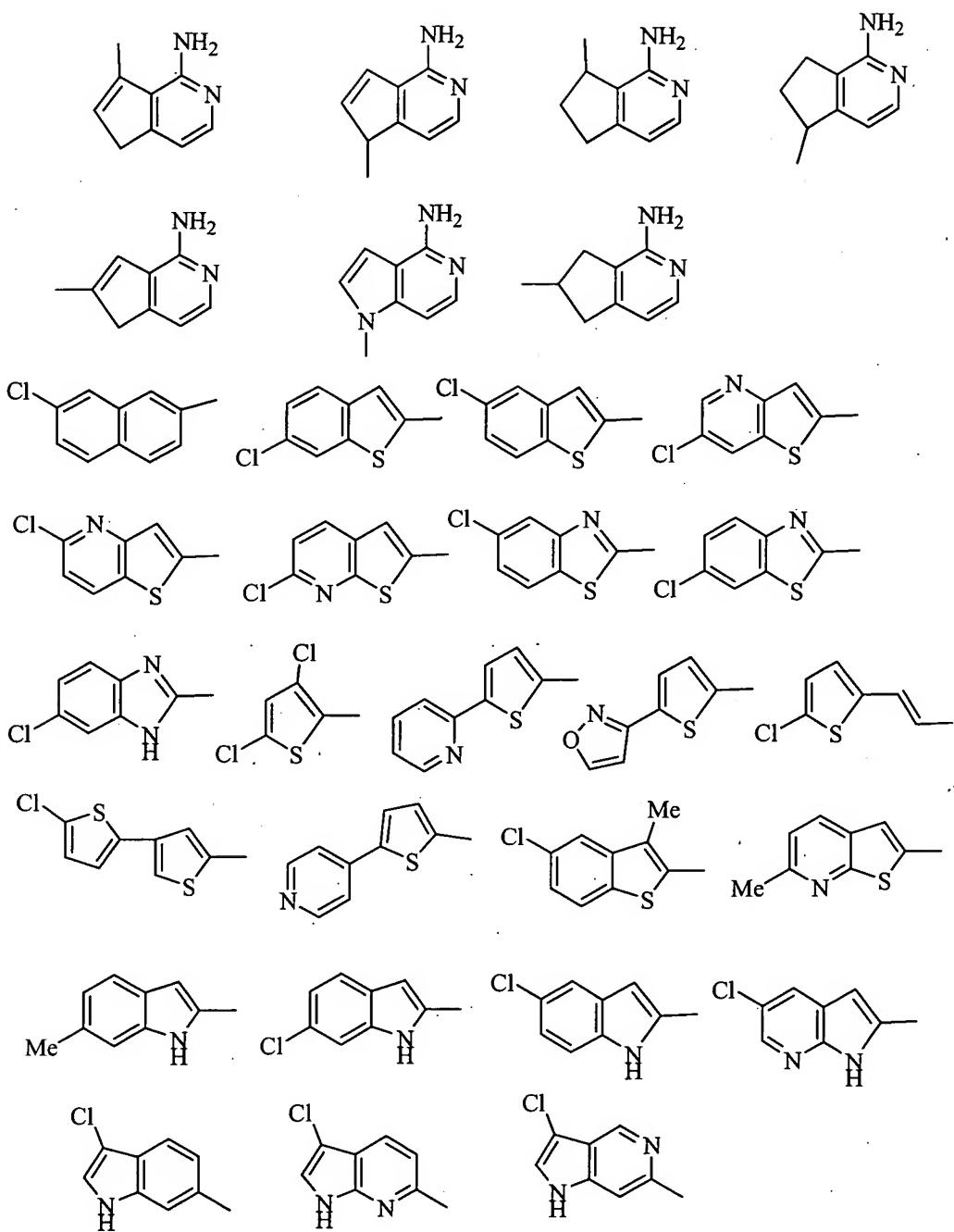


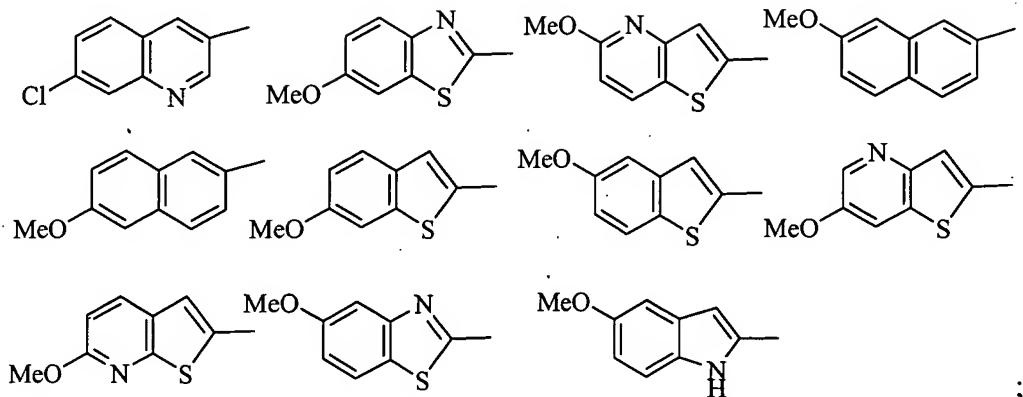






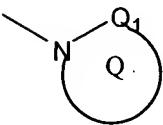






A is selected from one of the following carbocyclic and heterocyclic groups which are substituted with 0-2 R<sup>4</sup>:

cyclohexyl, phenyl, piperidinyl, piperazinyl, pyridyl, pyrimidyl, furanyl, morpholinyl, thienyl, pyrrolyl, pyrrolidinyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, pyrazolyl, imidazolyl, 1,2,3-oxadiazolyl, 1,2,4-oxadiazolyl, 1,2,5-oxadiazolyl, 1,3,4-oxadiazolyl, 1,2,3-thiadiazolyl, 1,2,4-thiadiazolyl, 1,2,5-thiadiazolyl, 1,3,4-thiadiazolyl, 1,2,3-triazolyl, 1,2,4-triazolyl, 1,2,5-triazolyl, 1,3,4-triazolyl, benzofuranyl, benzothiofuranyl, indolinyl, indolyl, benzimidazolyl, benzoxazolyl, benzthiazolyl, indazolyl, benzisoxazolyl, benzisothiazolyl, and isoindazolyl;

B is  ; provided that Z and B are attached to different atoms on A;

Q<sub>1</sub> is selected from C=O and SO<sub>2</sub>;

ring Q is a 5-7 membered ring consisting of, in addition to the amide group shown, carbon atoms and 0-2 heteroatoms selected from NR<sup>4c</sup>, O, S, S(O), and S(O)<sub>2</sub>, wherein:

0-2 double bonds are present within the ring and the ring is substituted with 0-2 R<sup>4a</sup>;

alternatively, ring Q is a 5-7 membered ring to which another ring is fused, wherein:

the 5-7 membered ring consists of, in addition to the shown amide group,

carbon atoms and 0-2 heteroatoms selected from NR<sup>4c</sup>, O, S, S(O), and S(O)<sub>2</sub>

and 0-1 double bonds are present within the ring;

the fusion ring is phenyl or a 5-6 membered heteroaromatic consisting of carbon atoms and 1-2 heteroatoms selected from NR<sup>4c</sup>, O, and S;

ring Q, which includes the 5-7 membered ring and the fusion ring, is substituted with 0-3 R<sup>4a</sup>;

R<sup>1a</sup>, at each occurrence, is selected from H, -(CH<sub>2</sub>)<sub>r</sub>-R<sup>1b</sup>, -(CH<sub>2</sub>)<sub>r</sub>-O-(CH<sub>2</sub>)<sub>r</sub>-R<sup>1b</sup>, -(CH<sub>2</sub>)<sub>r</sub>-C(=NR<sup>1b</sup>)NR<sup>3</sup>R<sup>1b</sup>, NR<sup>3</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>t</sub>R<sup>1c</sup>, O(CR<sup>3</sup>R<sup>3a</sup>)<sub>t</sub>R<sup>1c</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>(CH<sub>2</sub>)<sub>r</sub>R<sup>1b</sup>, (CH<sub>2</sub>)<sub>r</sub>C(O)NR<sup>2</sup>(CH<sub>2</sub>)<sub>r</sub>R<sup>1b</sup>, CO<sub>2</sub>(CH<sub>2</sub>)<sub>t</sub>R<sup>1b</sup>, O(CH<sub>2</sub>)<sub>t</sub>R<sup>1b</sup>, S(O)<sub>p</sub>(CH<sub>2</sub>)<sub>r</sub>R<sup>1d</sup>, O(CH<sub>2</sub>)<sub>r</sub>R<sup>1d</sup>, NR<sup>3</sup>(CH<sub>2</sub>)<sub>r</sub>R<sup>1d</sup>, OC(O)NR<sup>3</sup>(CH<sub>2</sub>)<sub>r</sub>R<sup>1d</sup>, NR<sup>3</sup>C(O)NR<sup>3</sup>(CH<sub>2</sub>)<sub>r</sub>R<sup>1d</sup>, NR<sup>3</sup>C(O)O(CH<sub>2</sub>)<sub>r</sub>R<sup>1d</sup>, and NR<sup>3</sup>C(O)(CH<sub>2</sub>)<sub>r</sub>R<sup>1d</sup>, provided that R<sup>1a</sup> forms other than an N-halo, N-S, O-O, or N-CN bond;

alternatively, when two R<sup>1a</sup> groups are attached to the same carbon atom, together with the carbon atom to which they are attached they form a 3-6 membered carbocyclic or heterocyclic ring consisting of: carbon atoms and 0-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub>, this ring being substituted with 0-2 R<sup>4</sup> and 0-3 ring double bonds;

R<sup>1b</sup> is selected from H, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, F, Cl, Br, -CN, -CHO, CF<sub>3</sub>, (CH<sub>2</sub>)<sub>r</sub>OR<sup>2</sup>, NR<sup>2</sup>R<sup>2a</sup>, C(O)R<sup>2b</sup>, CO<sub>2</sub>R<sup>2b</sup>, OC(O)R<sup>2</sup>, CO<sub>2</sub>R<sup>2a</sup>, S(O)<sub>p</sub>R<sup>2</sup>, NR<sup>2</sup>(CH<sub>2</sub>)<sub>r</sub>OR<sup>2</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>, NR<sup>2</sup>C(O)NR<sup>2</sup>R<sup>2a</sup>, C(O)NR<sup>2</sup>R<sup>2a</sup>, SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>R<sup>2</sup>, C(O)NR<sup>2</sup>SO<sub>2</sub>R<sup>2</sup>, SO<sub>2</sub>NR<sup>2</sup>C(O)R<sup>2</sup>, C<sub>3-10</sub> carbocycle substituted with 0-2 R<sup>4</sup>, and 4-10 membered heterocycle consisting of carbon atoms and

from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4b</sup>, provided that R<sup>1b</sup> forms other than an O-O, N-halo, N-S, or N-CN bond;

R<sup>2</sup>, at each occurrence, is selected from H, CF<sub>3</sub>, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, phenyl substituted with 0-2 R<sup>4b</sup>, a benzyl substituted with 0-2 R<sup>4b</sup>, and 5-6 membered aromatic heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4b</sup>;

R<sup>2a</sup>, at each occurrence, is selected from H, CF<sub>3</sub>, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, benzyl, phenyl substituted with 0-2 R<sup>4b</sup>, and 5-6 membered aromatic heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4b</sup>;

alternatively, R<sup>2</sup> and R<sup>2a</sup>, together with the atom to which they are attached, combine to form a 5 or 6 membered saturated, partially saturated or unsaturated ring substituted with 0-2 R<sup>4b</sup> and consisting of: 0-1 additional heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub>;

R<sup>2b</sup>, at each occurrence, is selected from CF<sub>3</sub>, C1-4 alkoxy, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, benzyl, phenyl substituted with 0-2 R<sup>4b</sup>, and 5-6 membered aromatic heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4b</sup>;

R<sup>2c</sup>, at each occurrence, is selected from CF<sub>3</sub>, OH, OCH<sub>3</sub>, OCH<sub>2</sub>CH<sub>3</sub>, OCH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, OCH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, benzyl, phenyl

substituted with 0-2 R<sup>4b</sup>, and 5-6 membered aromatic heterocycle containing from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4b</sup>;

R<sup>4</sup>, at each occurrence, is selected from H, =O, (CH<sub>2</sub>)<sub>r</sub>OR<sup>2</sup>, F, Cl, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, -CN, NO<sub>2</sub>, NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, C(O)R<sup>2c</sup>, CH<sub>2</sub>C(O)R<sup>2c</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>, CH<sub>2</sub>NR<sup>2</sup>C(O)R<sup>2b</sup>, C(O)NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>C(O)NR<sup>2</sup>R<sup>2a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>(CH<sub>2</sub>)<sub>1-2</sub>C(O)OR<sup>3</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>(CH<sub>2</sub>)<sub>2-4</sub>NR<sup>3</sup>R<sup>3a</sup>, (CH<sub>2</sub>)<sub>r</sub>NR<sup>3</sup>(CH<sub>2</sub>)<sub>2-4</sub>NR<sup>3</sup>SO<sub>2</sub>R<sup>3a</sup>, SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, CH<sub>2</sub>NR<sup>2</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, NR<sup>2</sup>SO<sub>2</sub>R<sup>5</sup>, CH<sub>2</sub>NR<sup>2</sup>SO<sub>2</sub>R<sup>5</sup>, S(O)<sub>p</sub>R<sup>5a</sup>, CH<sub>2</sub>S(O)<sub>p</sub>R<sup>5a</sup>, CF<sub>3</sub>, (CH<sub>2</sub>)<sub>r</sub>-3-7 membered carbocycle substituted with 0-1 R<sup>5</sup>, and a (CH<sub>2</sub>)<sub>r</sub>-5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-1 R<sup>5</sup>;

R<sup>4a</sup>, at each occurrence, is selected from H, =O, CH<sub>2</sub>OR<sup>2</sup>, OR<sup>2</sup>, F, Br, Cl, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH(CH<sub>3</sub>)<sub>2</sub>, CH(CH<sub>3</sub>)CH<sub>2</sub>CH<sub>3</sub>, C(CH<sub>3</sub>)<sub>3</sub>, -CN, NO<sub>2</sub>, CH<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>R<sup>2a</sup>, C(O)R<sup>2c</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>, C(O)NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>C(O)NR<sup>2</sup>R<sup>2a</sup>, SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, and -CF<sub>3</sub>;

R<sup>4b</sup>, at each occurrence, is selected from H, =O, (CH<sub>2</sub>)<sub>r</sub>OR<sup>3</sup>, F, Cl, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, -CN, NO<sub>2</sub>, NR<sup>3</sup>R<sup>3a</sup>, CH<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, C(O)R<sup>3</sup>, CH<sub>2</sub>C(O)R<sup>3</sup>, C(O)OR<sup>3c</sup>, CH<sub>2</sub>C(O)OR<sup>3c</sup>, NR<sup>3</sup>C(O)R<sup>3a</sup>, CH<sub>2</sub>NR<sup>3</sup>C(O)R<sup>3a</sup>, C(O)NR<sup>3</sup>R<sup>3a</sup>, CH<sub>2</sub>C(O)NR<sup>3</sup>R<sup>3a</sup>, SO<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, CH<sub>2</sub>SO<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, NR<sup>3</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, CH<sub>2</sub>NR<sup>3</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, NR<sup>3</sup>SO<sub>2</sub>-phenyl,

$\text{CH}_2\text{NR}^3\text{SO}_2\text{-phenyl}$ ,  $\text{S(O)pCF}_3$ ,  $\text{CH}_2\text{S(O)pCF}_3$ ,  $\text{S(O)p-C}_1\text{-4 alkyl}$ ,  
 $\text{CH}_2\text{S(O)p-C}_1\text{-4 alkyl}$ ,  $\text{S(O)p-phenyl}$ ,  $\text{CH}_2\text{S(O)p-phenyl}$ , and  $\text{CF}_3$ ;

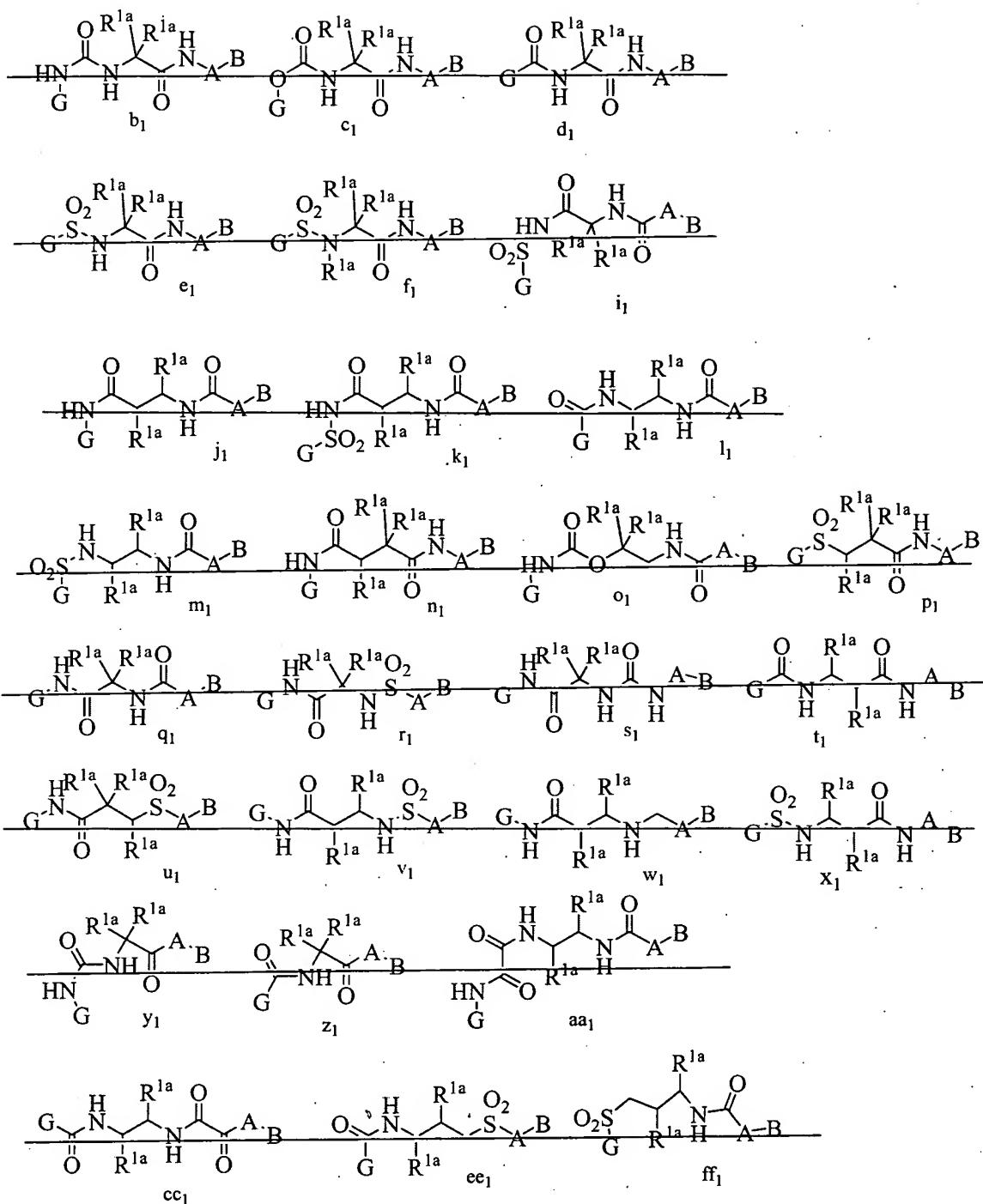
$\text{R}^{4c}$ , at each occurrence, is selected from  $\text{H}$ ,  $\text{CH}_3$ ,  $\text{CH}_2\text{CH}_3$ ,  $\text{CH}_2\text{CH}_2\text{CH}_3$ ,  $\text{CH}(\text{CH}_3)_2$ ,  
 $\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_3$ ,  $\text{CH}_2\text{CH}(\text{CH}_3)_2$ ,  $\text{CH}(\text{CH}_3)\text{CH}_2\text{CH}_3$ ,  $\text{C}(\text{CH}_3)_3$ ,  $\text{CH}_2\text{OR}^2$ ,  
 $\text{CH}_2\text{F}$ ,  $\text{CH}_2\text{Br}$ ,  $\text{CH}_2\text{Cl}$ ,  $\text{CH}_2\text{CN}$ ,  $\text{CH}_2\text{NO}_2$ ,  $\text{CH}_2\text{NR}^2\text{R}^{2a}$ ,  $\text{C(O)R}^{2c}$ ,  
 $\text{CH}_2\text{C(O)R}^{2c}$ ,  $\text{CH}_2\text{NR}^2\text{C(O)R}^{2b}$ ,  $\text{C(O)NR}^2\text{R}^{2a}$ ,  $\text{CH}_2\text{C(O)NR}^2\text{R}^{2a}$ ,  
 $\text{SO}_2\text{NR}^2\text{R}^{2a}$ ,  $\text{CH}_2\text{SO}_2\text{NR}^2\text{R}^{2a}$ ,  $\text{S(O)pR}^{5a}$ ,  $\text{CH}_2\text{S(O)pR}^{5a}$ ,  $\text{CF}_3$ , phenyl  
substituted with 0-1  $\text{R}^5$ , and benzyl substituted with 0-1  $\text{R}^5$ ;

$\text{R}^5$ , at each occurrence, is selected from  $\text{H}$ ,  $=\text{O}$ ,  $\text{CH}_3$ ,  $\text{CH}_2\text{CH}_3$ ,  $\text{CH}_2\text{CH}_2\text{CH}_3$ ,  
 $\text{CH}(\text{CH}_3)_2$ ,  $\text{OR}^3$ ,  $\text{CH}_2\text{OR}^3$ ,  $\text{F}$ ,  $\text{Cl}$ ,  $-\text{CN}$ ,  $\text{NO}_2$ ,  $\text{NR}^3\text{R}^{3a}$ ,  $\text{CH}_2\text{NR}^3\text{R}^{3a}$ ,  $\text{C(O)R}^3$ ,  
 $\text{CH}_2\text{C(O)R}^3$ ,  $\text{C(O)OR}^{3c}$ ,  $\text{CH}_2\text{C(O)OR}^{3c}$ ,  $\text{NR}^3\text{C(O)R}^{3a}$ ,  $\text{C(O)NR}^3\text{R}^{3a}$ ,  
 $\text{SO}_2\text{NR}^3\text{R}^{3a}$ ,  $\text{NR}^3\text{SO}_2\text{-C}_1\text{-4 alkyl}$ ,  $\text{NR}^3\text{SO}_2\text{CF}_3$ ,  $\text{NR}^3\text{SO}_2\text{-phenyl}$ ,  $\text{S(O)pCF}_3$ ,  
 $\text{S(O)p-C}_1\text{-4 alkyl}$ ,  $\text{S(O)p-phenyl}$ ,  $\text{CF}_3$ , phenyl substituted with 0-2  $\text{R}^6$ , naphthyl  
substituted with 0-2  $\text{R}^6$ , and benzyl substituted with 0-2  $\text{R}^6$ ;

$\text{R}^6$ , at each occurrence, is selected from  $\text{H}$ ,  $\text{OH}$ ,  $\text{OR}^2$ ,  $\text{F}$ ,  $\text{Cl}$ ,  $\text{CH}_3$ ,  $\text{CH}_2\text{CH}_3$ ,  
 $\text{CH}_2\text{CH}_2\text{CH}_3$ ,  $\text{CH}(\text{CH}_3)_2$ ,  $-\text{CN}$ ,  $\text{NO}_2$ ,  $\text{NR}^2\text{R}^{2a}$ ,  $\text{CH}_2\text{NR}^2\text{R}^{2a}$ ,  $\text{C(O)R}^{2b}$ ,  
 $\text{CH}_2\text{C(O)R}^{2b}$ ,  $\text{NR}^2\text{C(O)R}^{2b}$ ,  $\text{SO}_2\text{NR}^2\text{R}^{2a}$ , and  $\text{NR}^2\text{SO}_2\text{C}_1\text{-4 alkyl}$ ; and,

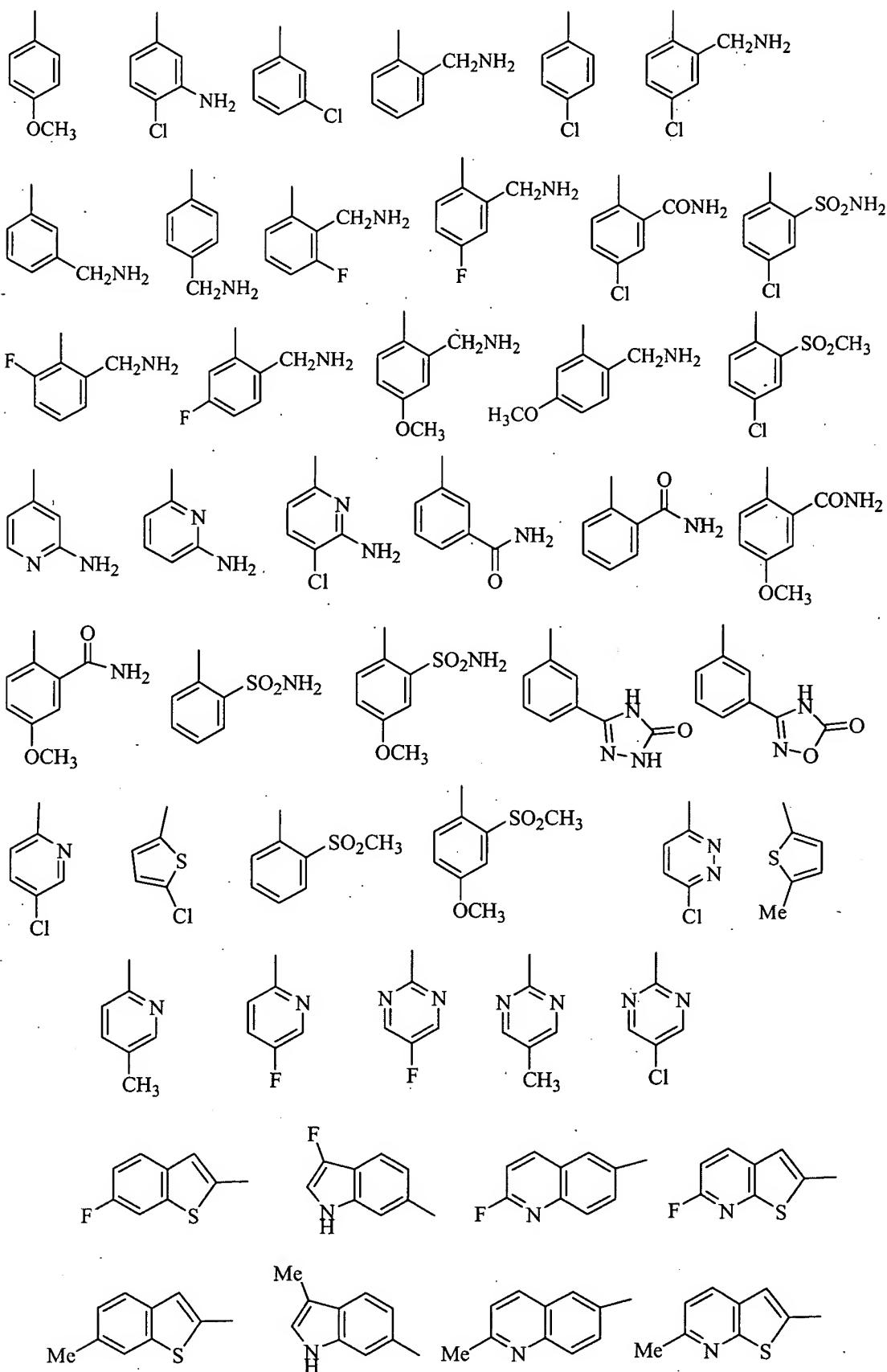
$r$ , at each occurrence, is selected from 0, 1, and 2.

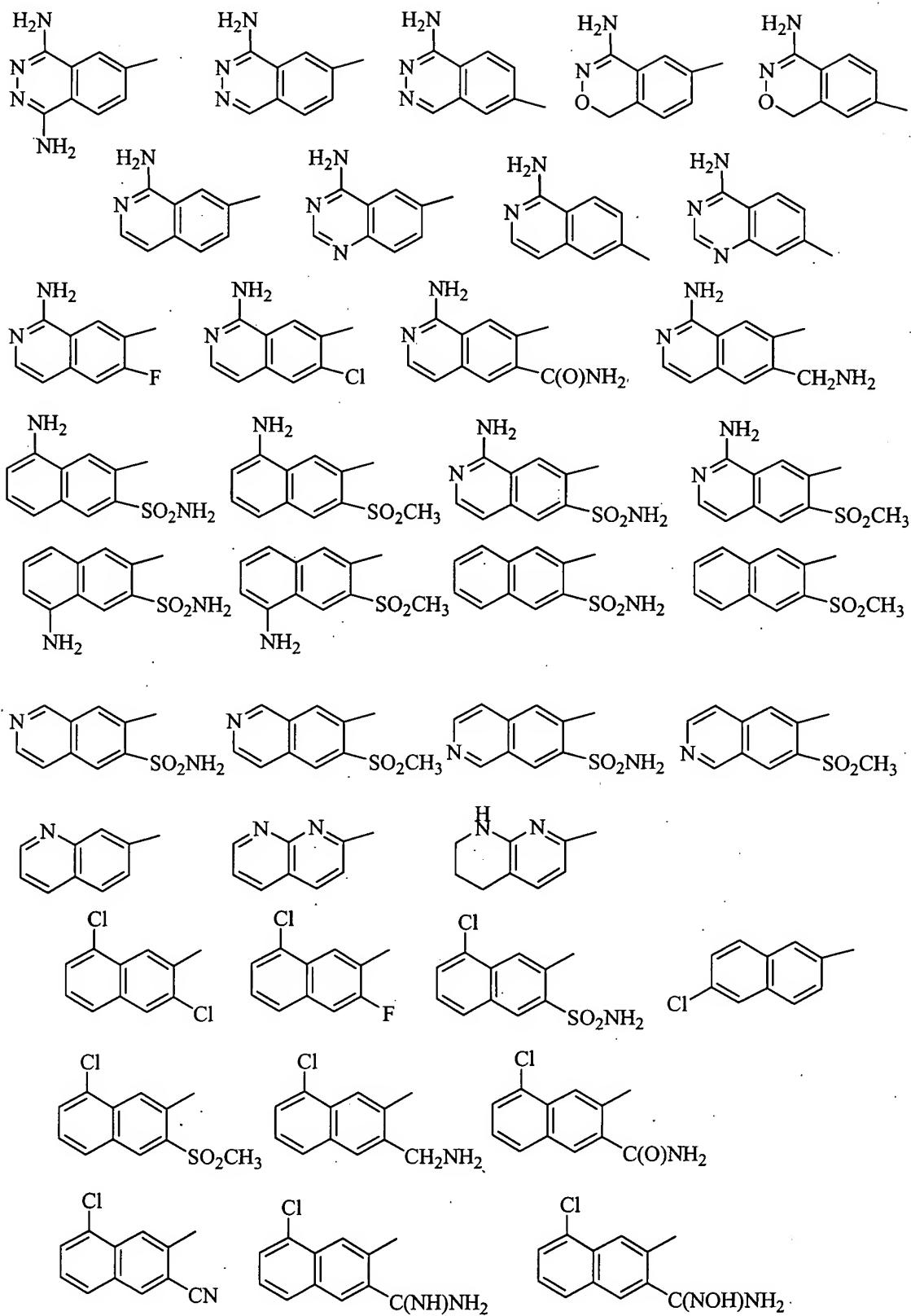
4. (Currently Amended) A compound according to Claim 3, wherein ~~the compound is selected from compounds b<sub>1</sub>-f<sub>1</sub>, i<sub>1</sub>-aa<sub>1</sub>, ee<sub>1</sub>, ee<sub>1</sub>, and ff<sub>1</sub>~~:

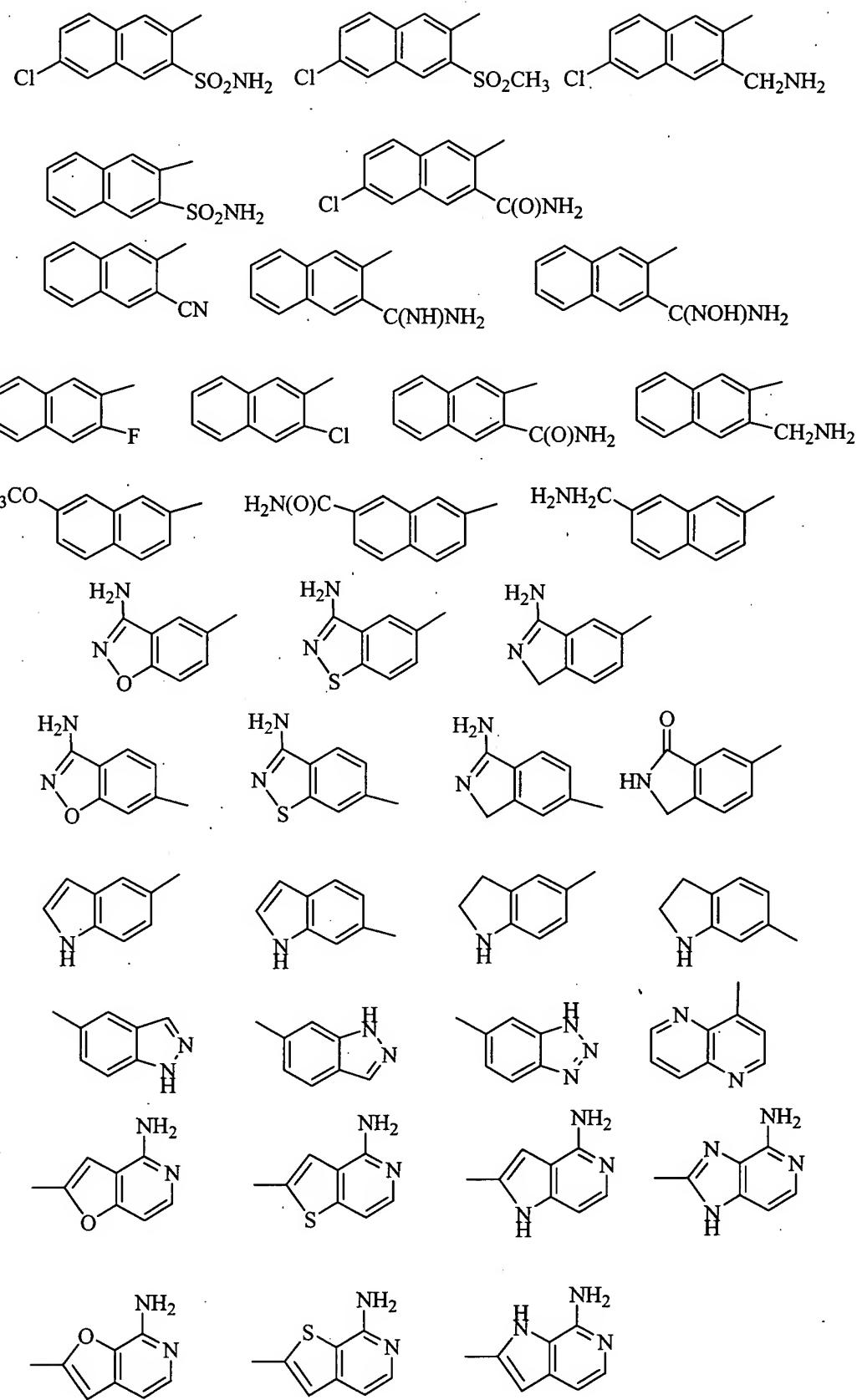


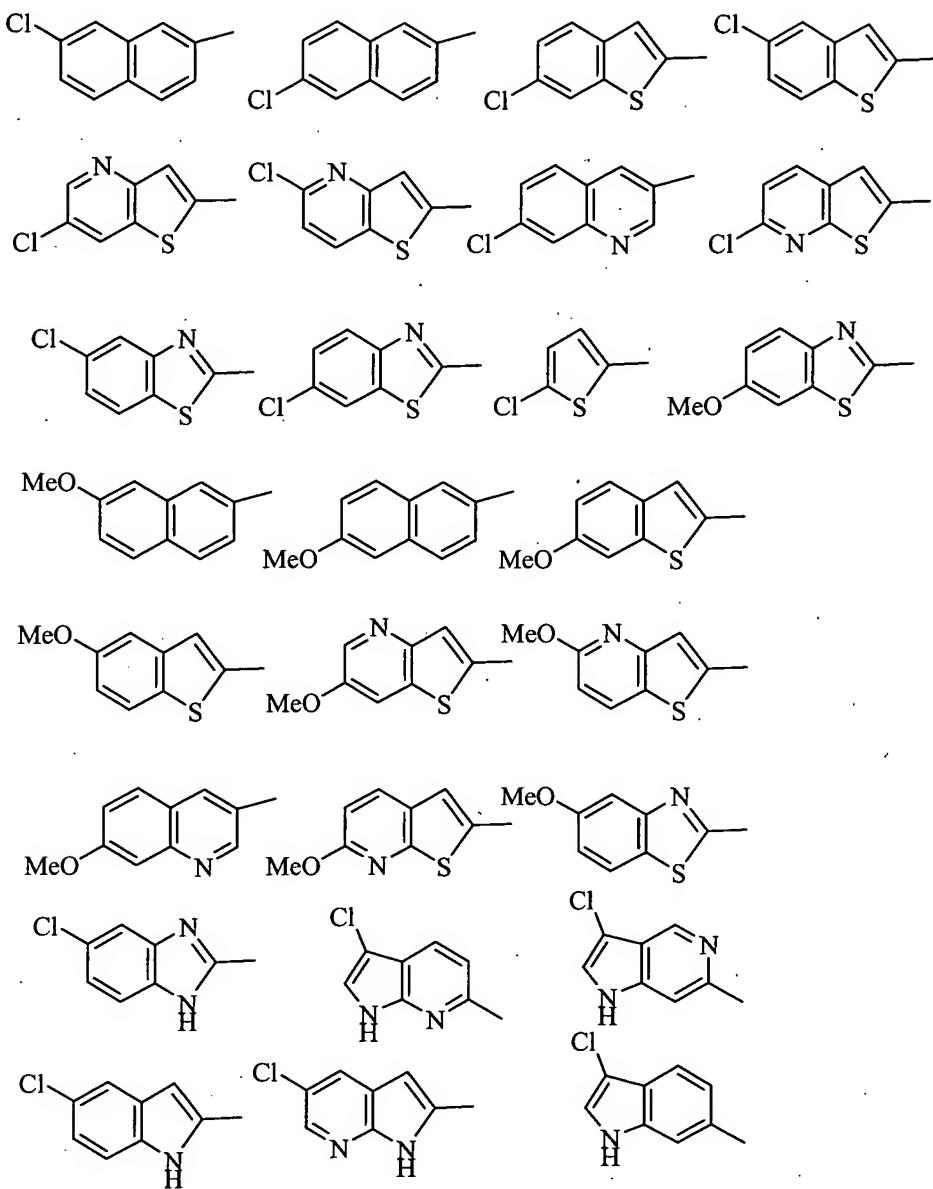
**wherein:**

$G$  is selected from the group:

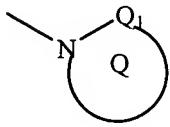








A is selected from cyclohexyl, piperidinyl, indolinyl, phenyl, pyridyl, thienyl, and pyrimidyl, and is substituted with 0-2 R<sup>4</sup>;



B is ; provided that Z and B are attached to different atoms on A;

Q<sub>1</sub> is selected from C=O and SO<sub>2</sub>;

ring Q is a 5-6 membered ring consisting of, in addition to the amide group shown, carbon atoms and 0-1 heteroatoms selected from NR<sup>4c</sup>, O, S, S(O), and S(O)<sub>2</sub>, wherein:

0-2 double bonds are present within the ring and the ring is substituted with 0-2 R<sup>4a</sup>;

alternatively, ring Q is a 5-7 membered ring to which another ring is fused, wherein:

the 5-7 membered ring consists of, in addition to the shown amide group, carbon atoms and 0-1 heteroatoms selected from NR<sup>4c</sup>, O, S, S(O), and S(O)<sub>2</sub> } and 0-1 double bonds are present within the ring;

the fusion ring is phenyl;  
ring Q, which includes the 5-7 membered ring and the fusion ring, is substituted with 0-2R<sup>4a</sup>;

R<sup>1a</sup> is selected from H, R<sup>1b</sup>, C(CH<sub>3</sub>)<sub>2</sub>R<sup>1b</sup>, CH(CH<sub>3</sub>)R<sup>1b</sup>, CH<sub>2</sub>R<sup>1b</sup>, CH<sub>2</sub>CH<sub>2</sub>R<sup>1b</sup>, CH<sub>2</sub>OCH<sub>2</sub>CH<sub>2</sub>R<sup>1b</sup>, OCH<sub>2</sub>CH<sub>2</sub>R<sup>1b</sup>, (CH<sub>2</sub>)<sub>t</sub>NR<sup>3</sup>CH<sub>2</sub>CH<sub>2</sub>R<sup>1b</sup>, NR<sup>3</sup>(CR<sup>3</sup>R<sup>3a</sup>)<sub>t</sub>R<sup>1c</sup>, O(CR<sup>3</sup>R<sup>3a</sup>)<sub>t</sub>R<sup>1c</sup>, (CH<sub>2</sub>)<sub>t</sub>C(O)NR<sup>2</sup>(CH<sub>2</sub>)<sub>t</sub>R<sup>1b</sup>, S(O)<sub>p</sub>(CH<sub>2</sub>)<sub>t</sub>R<sup>1d</sup>, O(CH<sub>2</sub>)<sub>t</sub>R<sup>1d</sup>, NR<sup>3</sup>(CH<sub>2</sub>)<sub>t</sub>R<sup>1d</sup>, OC(O)NR<sup>3</sup>(CH<sub>2</sub>)<sub>t</sub>R<sup>1d</sup>, NR<sup>3</sup>C(O)NR<sup>3</sup>(CH<sub>2</sub>)<sub>t</sub>R<sup>1d</sup>, NR<sup>3</sup>C(O)O(CH<sub>2</sub>)<sub>t</sub>R<sup>1d</sup>, and NR<sup>3</sup>C(O)(CH<sub>2</sub>)<sub>t</sub>R<sup>1d</sup>, provided that R<sup>1a</sup> forms other than an N-halo, N-S, O-O, or N-CN bond;

alternatively, when two R<sup>1a</sup> groups are attached to the same carbon atom, together with the carbon atom to which they are attached they form a 3-10 membered carbocyclic or heterocyclic ring consisting of: carbon atoms and 0-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub>, this ring being substituted with 0-2 R<sup>4</sup> and 0-2 ring double bonds;

$R^{1b}$  is selected from H, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, F, Cl, Br, -CN, -CHO, CF<sub>3</sub>, (CH<sub>2</sub>)<sub>r</sub>OR<sup>2</sup>, NR<sup>2</sup>R<sup>2a</sup>, C(O)R<sup>2b</sup>, CO<sub>2</sub>R<sup>2b</sup>, OC(O)R<sup>2</sup>, CO<sub>2</sub>R<sup>2a</sup>, S(O)<sub>p</sub>R<sup>2</sup>, NR<sup>2</sup>(CH<sub>2</sub>)<sub>r</sub>OR<sup>2</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>, NR<sup>2</sup>C(O)NR<sup>2</sup>R<sup>2a</sup>, C(O)NR<sup>2</sup>R<sup>2a</sup>, SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>R<sup>2</sup>, C(O)NR<sup>2</sup>SO<sub>2</sub>R<sup>2</sup>, SO<sub>2</sub>NR<sup>2</sup>C(O)R<sup>2</sup>, C<sub>3-6</sub> carbocycle substituted with 0-2 R<sup>4</sup>, and 4-10 membered heterocycle consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4</sup>, provided that R<sup>1b</sup> forms other than an O-O, N-halo, N-S, or N-CN bond;

R<sup>2</sup>, at each occurrence, is selected from H, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, phenyl substituted with 0-1 R<sup>4b</sup>, benzyl substituted with 0-1 R<sup>4b</sup>, and 5-6 membered aromatic heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-1 R<sup>4b</sup>;

R<sup>2a</sup>, at each occurrence, is selected from H, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, benzyl, phenyl substituted with 0-1 R<sup>4b</sup>, and 5-6 membered aromatic heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-1 R<sup>4b</sup>;

alternatively, R<sup>2</sup> and R<sup>2a</sup>, together with the atom to which they are attached, combine to form a 5 or 6 membered saturated, partially saturated or unsaturated ring substituted with 0-1 R<sup>4b</sup> and consisting of: 0-1 additional heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub>;

R<sup>2b</sup>, at each occurrence, is selected from OCH<sub>3</sub>, OCH<sub>2</sub>CH<sub>3</sub>, OCH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, OCH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, benzyl, phenyl

substituted with 0-1 R<sup>4b</sup>, and 5-6 membered aromatic heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-1 R<sup>4b</sup>;

R<sup>2c</sup>, at each occurrence, is selected from OH, OCH<sub>3</sub>, OCH<sub>2</sub>CH<sub>3</sub>, OCH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, OCH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, benzyl, phenyl substituted with 0-1 R<sup>4b</sup>, and 5-6 membered aromatic heterocycle containing from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-1 R<sup>4b</sup>;

R<sup>4</sup>, at each occurrence, is selected from H, =O, OR<sup>2</sup>, CH<sub>2</sub>OR<sup>2</sup>, F, Cl, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, -CN, NO<sub>2</sub>, NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, C(O)R<sup>2c</sup>, CH<sub>2</sub>C(O)R<sup>2c</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>, C(O)NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>C(O)NR<sup>2</sup>R<sup>2a</sup>, NR<sup>3</sup>(CH<sub>2</sub>)<sub>1-2</sub>C(O)OR<sup>3</sup>, NR<sup>3</sup>(CH<sub>2</sub>)<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, NR<sup>3</sup>(CH<sub>2</sub>)<sub>2</sub>NR<sup>3</sup>C(O)R<sup>3a</sup>, NR<sup>3</sup>(CH<sub>2</sub>)<sub>2</sub>NR<sup>3</sup>SO<sub>2</sub>R<sup>3a</sup>, SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, NR<sup>2</sup>SO<sub>2</sub>R<sup>5</sup>, S(O)<sub>p</sub>R<sup>5a</sup>, CF<sub>3</sub>, (CH<sub>2</sub>)<sub>1-3</sub> membered carbocycle substituted with 0-1 R<sup>5</sup>, and a (CH<sub>2</sub>)<sub>1-5</sub>-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-1 R<sup>5</sup>;

R<sup>4a</sup>, at each occurrence, is selected from H, =O, CH<sub>2</sub>OR<sup>2</sup>, OR<sup>2</sup>, F, Br, Cl, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH(CH<sub>3</sub>)<sub>2</sub>, CH(CH<sub>3</sub>)CH<sub>2</sub>CH<sub>3</sub>, C(CH<sub>3</sub>)<sub>3</sub>, CH<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, NR<sup>2</sup>R<sup>2a</sup>, C(O)R<sup>2c</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>, C(O)NR<sup>2</sup>R<sup>2a</sup>, SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, and CF<sub>3</sub>;

R<sup>4b</sup>, at each occurrence, is selected from H, =O, OR<sup>3</sup>, CH<sub>2</sub>OR<sup>3</sup>, F, Cl, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, -CN, NO<sub>2</sub>, NR<sup>3</sup>R<sup>3a</sup>, CH<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, C(O)R<sup>3</sup>,

C(O)OR<sup>3c</sup>, CH<sub>2</sub>C(O)OR<sup>3c</sup>, NR<sup>3</sup>C(O)R<sup>3a</sup>, C(O)NR<sup>3</sup>R<sup>3a</sup>, CH<sub>2</sub>C(O)NR<sup>3</sup>R<sup>3a</sup>, SO<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, NR<sup>3</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, NR<sup>3</sup>SO<sub>2</sub>-phenyl, S(O)p-C<sub>1-4</sub> alkyl, S(O)p-phenyl, and CF<sub>3</sub>;

R<sup>4c</sup>, at each occurrence, is selected from H, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, phenyl substituted with 0-1 R<sup>5</sup>, and benzyl substituted with 0-1 R<sup>5</sup>;

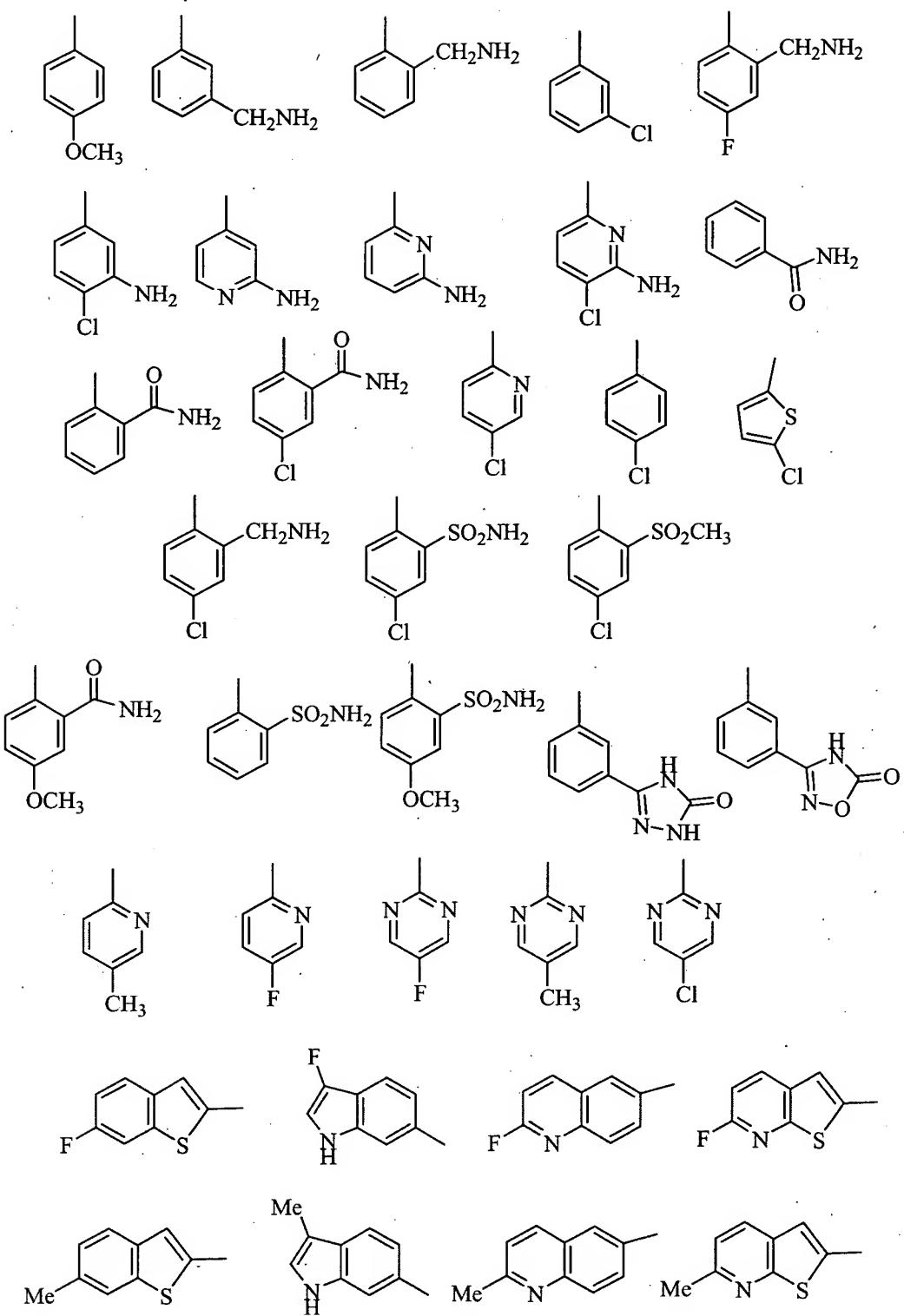
R<sup>5</sup>, at each occurrence, is selected from H, =O, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, OR<sup>3</sup>, CH<sub>2</sub>OR<sup>3</sup>, F, Cl, -CN, NO<sub>2</sub>, NR<sup>3</sup>R<sup>3a</sup>, CH<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, C(O)R<sup>3</sup>, C(O)OR<sup>3c</sup>, NR<sup>3</sup>C(O)R<sup>3a</sup>, C(O)NR<sup>3</sup>R<sup>3a</sup>, SO<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, NR<sup>3</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, NR<sup>3</sup>SO<sub>2</sub>-phenyl, S(O)p-C<sub>1-4</sub> alkyl, S(O)p-phenyl, CF<sub>3</sub>, phenyl substituted with 0-2 R<sup>6</sup>, naphthyl substituted with 0-2 R<sup>6</sup>, and benzyl substituted with 0-2 R<sup>6</sup>; and,

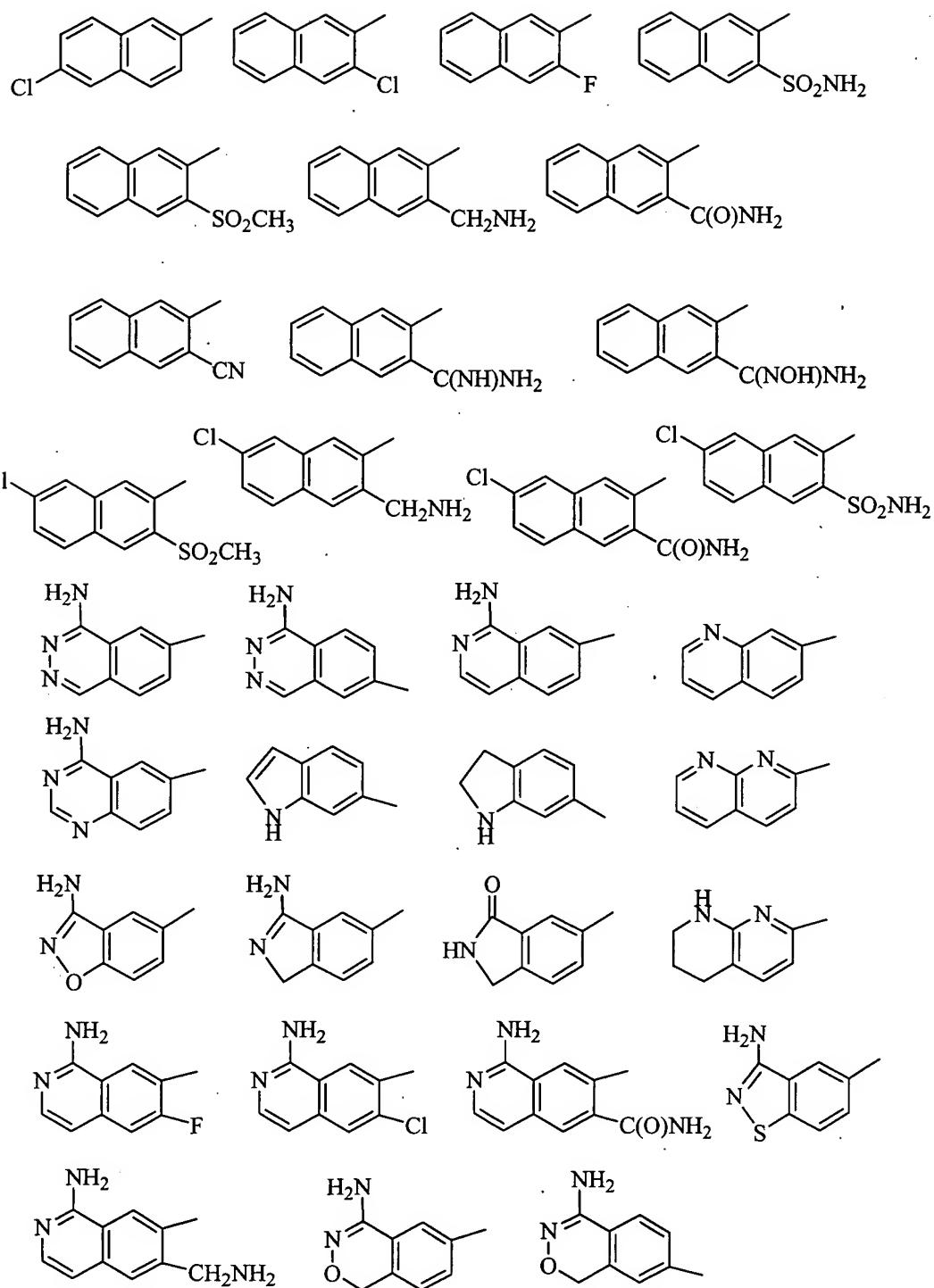
R<sup>6</sup>, at each occurrence, is selected from H, OH, OR<sup>2</sup>, F, Cl, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, -CN, NO<sub>2</sub>, NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, C(O)R<sup>2b</sup>, CH<sub>2</sub>C(O)R<sup>2b</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>, and SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>.

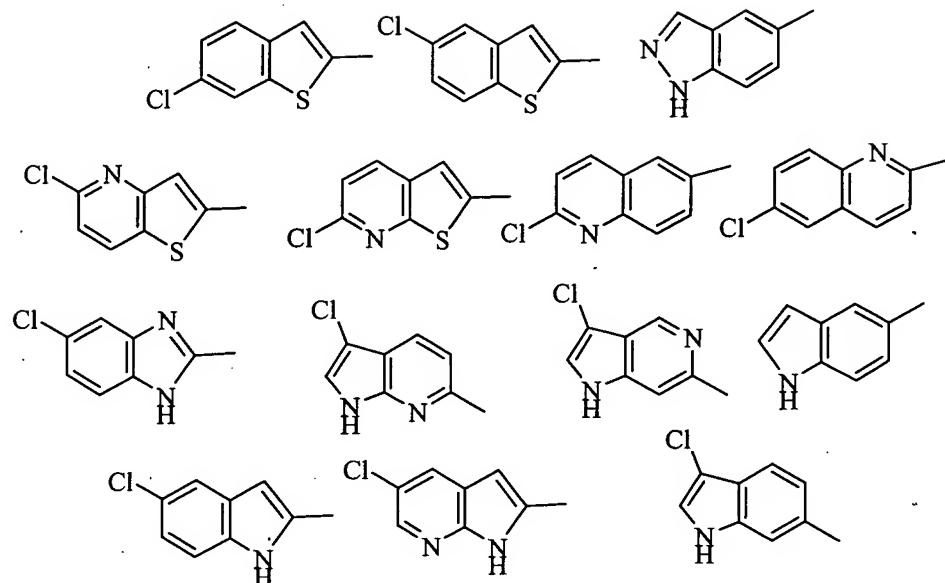
5. (Currently Amended) A compound according to Claim 4, wherein:

~~M is 4-7 membered linear chain consisting of: carbon atoms, 1-2 carbonyl groups, and 1-3 heteroatoms selected from O, S(O)p, and N, and M is substituted with 0-3 R<sup>1a</sup> and 0-1 R<sup>2</sup>, provided that other than an S-S, S-O, or O-O bond is present in M;~~

G is selected from:

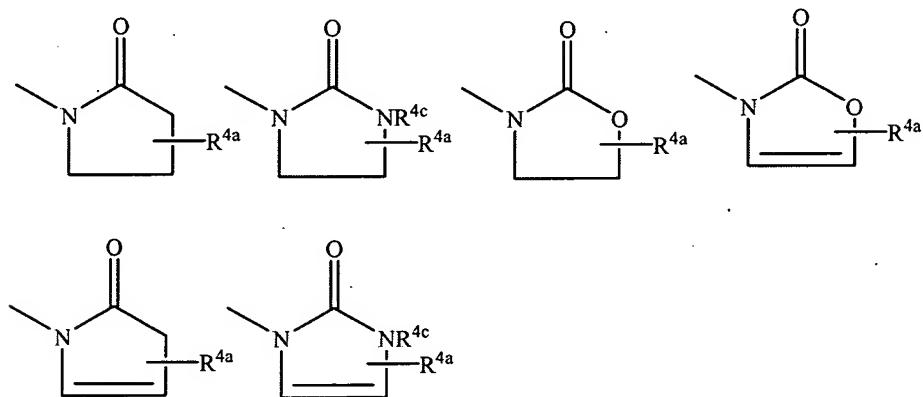


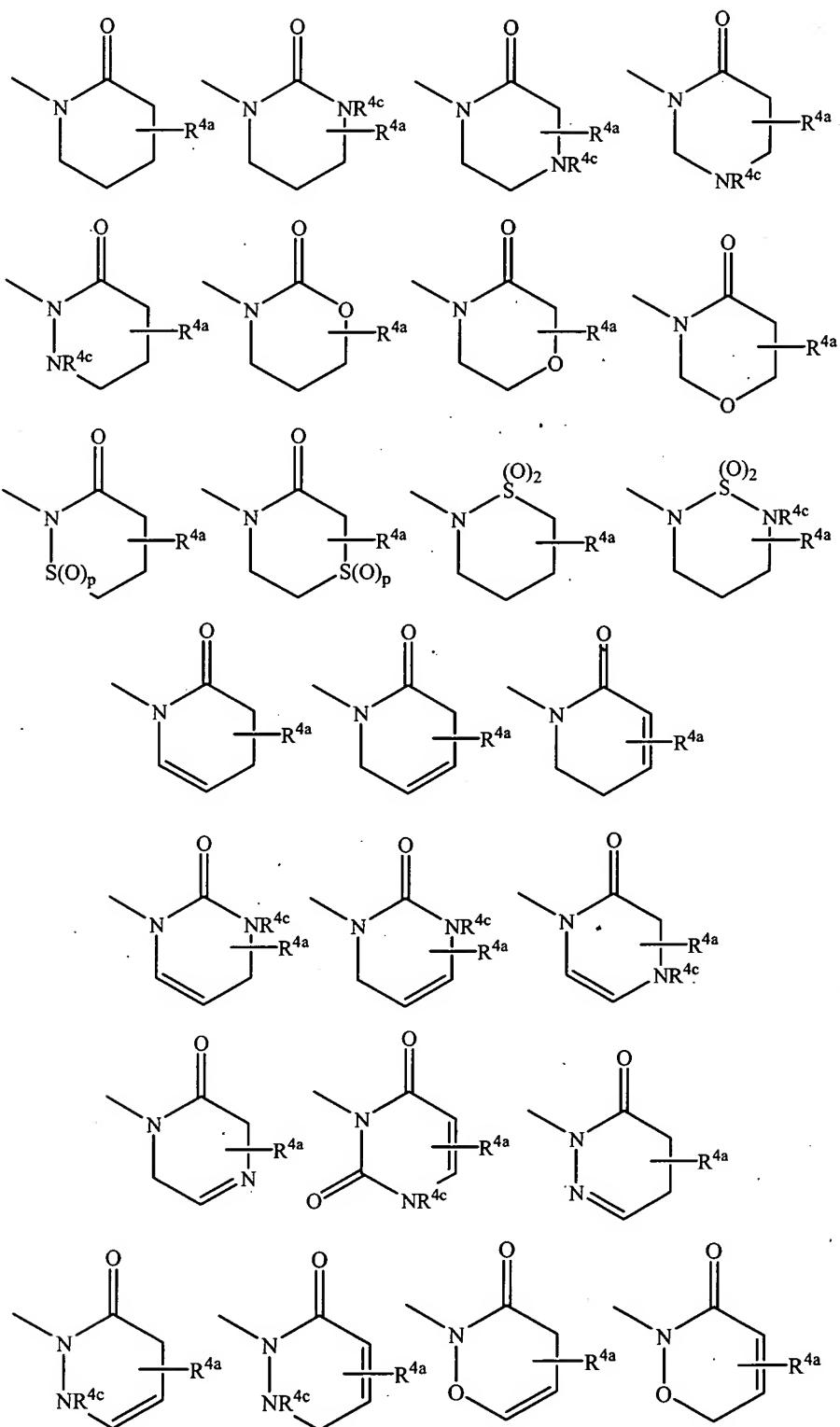


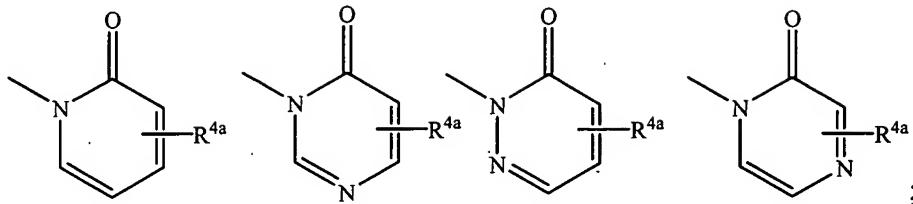


A is selected from the group: cyclohexyl, piperidinyl, indolinyl, phenyl, 2-pyridyl, 3-pyridyl, 2-pyrimidyl, 2-Cl-phenyl, 3-Cl-phenyl, 2-F-phenyl, 3-F-phenyl, 2-methylphenyl, 2-aminophenyl, and 2-methoxyphenyl;

B is attached to a different atom on A than Z and is selected from the group:







R<sup>1a</sup> is selected from H, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>(CH<sub>3</sub>)<sub>2</sub>, CF<sub>3</sub>, CH<sub>2</sub>CF<sub>3</sub>, OCH<sub>3</sub>, CH<sub>2</sub>OH, C(CH<sub>3</sub>)<sub>2</sub>OH, CH<sub>2</sub>OCH<sub>3</sub>, NH<sub>2</sub>, CH<sub>2</sub>NH<sub>2</sub>, NHCH<sub>3</sub>, CH<sub>2</sub>NHCH<sub>3</sub>, N(CH<sub>3</sub>)<sub>2</sub>, CH<sub>2</sub>N(CH<sub>3</sub>)<sub>2</sub>, CO<sub>2</sub>H, COCH<sub>3</sub>, CO<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CO<sub>2</sub>CH<sub>3</sub>, NHCOCH<sub>3</sub>, S(O)CH<sub>3</sub>, CH<sub>2</sub>S(O)CH<sub>3</sub>, S(O)<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>S(O)<sub>2</sub>CH<sub>3</sub>, C(O)NH<sub>2</sub>, CH<sub>2</sub>C(O)NH<sub>2</sub>, SO<sub>2</sub>NH<sub>2</sub>, CH<sub>2</sub>SO<sub>2</sub>NH<sub>2</sub>, NHSO<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>NHSO<sub>2</sub>CH<sub>3</sub>, NHSO<sub>2</sub>NHCH<sub>3</sub>, NHSO<sub>2</sub>N(CH<sub>3</sub>)<sub>2</sub>, NHCO<sub>2</sub>R<sup>2a</sup>, NHC(O)NHR<sup>2a</sup>, CH<sub>2</sub>OCH<sub>2</sub>CH<sub>2</sub>NR<sup>2a</sup>, C(O)NR<sup>2a</sup>, CH<sub>2</sub>CH<sub>2</sub>OR<sup>2</sup>, CH<sub>2</sub>C(O)NR<sup>2a</sup>CH<sub>2</sub>CH<sub>2</sub>OR<sup>2</sup>, C(O)NHCH<sub>2</sub>CH<sub>2</sub>NR<sup>2a</sup>, CH<sub>2</sub>C(O)NCH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>NR<sup>2a</sup>, CH<sub>2</sub>C(O)NCH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>NR<sup>2a</sup>, CH<sub>2</sub>N(CH<sub>3</sub>)CH<sub>2</sub>CH<sub>2</sub>NR<sup>2a</sup>, phenyl substituted with 0-2 R<sup>4b</sup>, -CH<sub>2</sub>-phenyl substituted with 0-2 R<sup>4b</sup>, 5-10 membered aromatic heterocycle consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4b</sup>, and -CH<sub>2</sub>-5-10 membered aromatic heterocycle consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-2 R<sup>4b</sup>, provided that R<sup>1a</sup> forms other than an N-halo, N-S, O-O, or N-CN bond;

R<sup>2</sup>, at each occurrence, is selected from H, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, phenyl substituted with 0-1 R<sup>4b</sup>, benzyl substituted with 0-1 R<sup>4b</sup>, and 5 membered aromatic heterocycle consisting of: carbon atoms and 1-4 heteroatoms

selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-1 R<sup>4b</sup>;

R<sup>2a</sup>, at each occurrence, is selected from H, CH<sub>3</sub>, and CH<sub>2</sub>CH<sub>3</sub>;

alternatively, R<sup>2</sup> and R<sup>2a</sup>, together with the atom to which they are attached, combine to form a 5 or 6 membered saturated, partially saturated or unsaturated ring substituted with 0-1 R<sup>4b</sup> and consisting of: 0-1 additional heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub>;

R<sup>2b</sup>, at each occurrence, is selected from OH, OCH<sub>3</sub>, OCH<sub>2</sub>CH<sub>3</sub>, CH<sub>3</sub>, and CH<sub>2</sub>CH<sub>3</sub>;

R<sup>2c</sup>, at each occurrence, is selected from OH, OCH<sub>3</sub>, OCH<sub>2</sub>CH<sub>3</sub>, CH<sub>3</sub>, and CH<sub>2</sub>CH<sub>3</sub>;

R<sup>4</sup>, at each occurrence, is selected from H, =O, OR<sup>2</sup>, CH<sub>2</sub>OR<sup>2</sup>, F, Cl, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, C(O)R<sup>2c</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>, C(O)NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>C(O)NR<sup>2</sup>R<sup>2a</sup>, NR<sup>3</sup>CH<sub>2</sub>C(O)OR<sup>3</sup>, NR<sup>3</sup>CH<sub>2</sub>CH<sub>2</sub>C(O)OR<sup>3</sup>, NR<sup>3</sup>(CH<sub>2</sub>)<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, NR<sup>3</sup>(CH<sub>2</sub>)<sub>2</sub>NR<sup>3</sup>C(O)R<sup>3a</sup>, NR<sup>3</sup>(CH<sub>2</sub>)<sub>2</sub>NR<sup>3</sup>SO<sub>2</sub>R<sup>3a</sup>, NR<sup>2</sup>SO<sub>2</sub>R<sup>5</sup>, S(O)<sub>2</sub>CH<sub>3</sub>, S(O)<sub>2</sub>-phenyl, CF<sub>3</sub>, (CH<sub>2</sub>)<sub>r</sub>-3-7 membered carbocycle substituted with 0-1 R<sup>5</sup>, and a (CH<sub>2</sub>)<sub>r</sub>-5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)<sub>p</sub> and substituted with 0-1 R<sup>5</sup>;

R<sup>4a</sup>, at each occurrence, is selected from H, =O, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH(CH<sub>3</sub>)<sub>2</sub>, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>, CH<sub>2</sub>CH(CH<sub>3</sub>)<sub>2</sub>, CH(CH<sub>3</sub>)CH<sub>2</sub>CH<sub>3</sub>, and C(CH<sub>3</sub>)<sub>3</sub>;

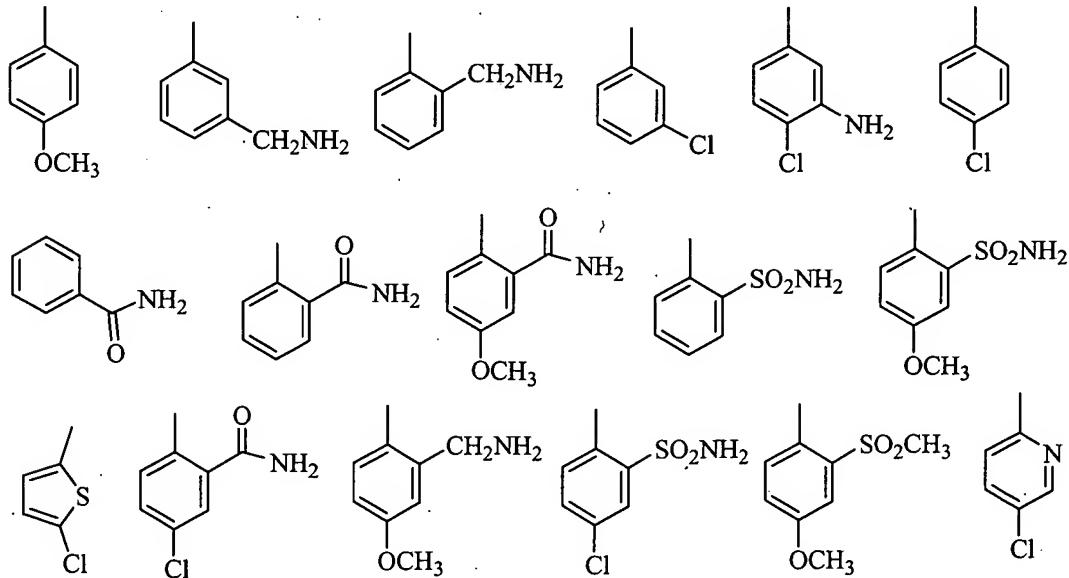
$R^{4b}$ , at each occurrence, is selected from H, =O, OR<sup>3</sup>, CH<sub>2</sub>OR<sup>3</sup>, F, Cl, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, NR<sup>3</sup>R<sup>3a</sup>, CH<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, C(O)R<sup>3</sup>, C(O)OR<sup>3c</sup>, NR<sup>3</sup>C(O)R<sup>3a</sup>, C(O)NR<sup>3</sup>R<sup>3a</sup>, CH<sub>2</sub>C(O)NR<sup>3</sup>R<sup>3a</sup>, NR<sup>3</sup>SO<sub>2</sub>-phenyl, S(O)<sub>2</sub>CH<sub>3</sub>, S(O)<sub>2</sub>-phenyl, and CF<sub>3</sub>;

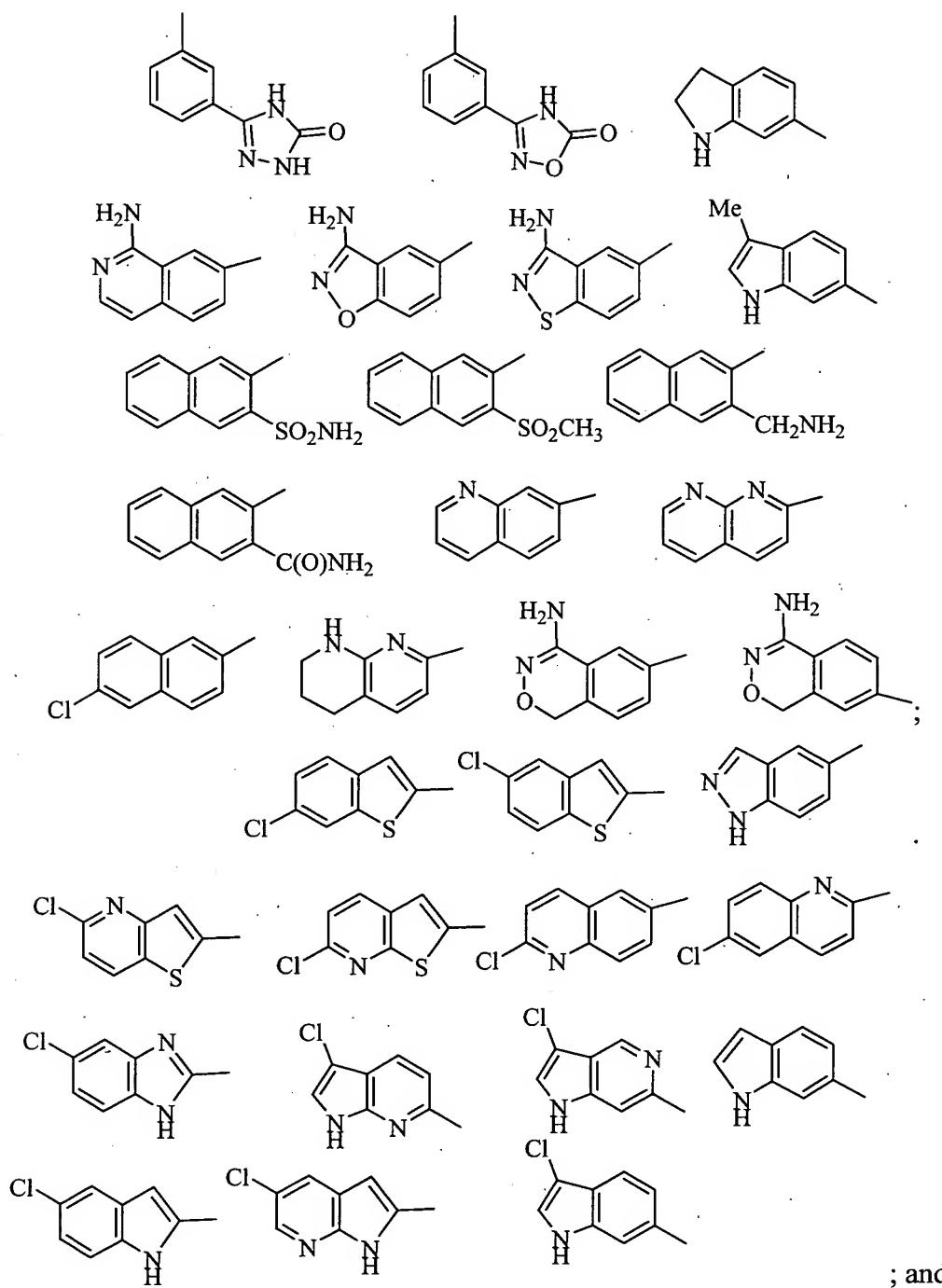
$R^5$ , at each occurrence, is selected from H, =O, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, OR<sup>3</sup>, CH<sub>2</sub>OR<sup>3</sup>, F, Cl, NR<sup>3</sup>R<sup>3a</sup>, CH<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, C(O)R<sup>3</sup>, C(O)OR<sup>3c</sup>, NR<sup>3</sup>C(O)R<sup>3a</sup>, C(O)NR<sup>3</sup>R<sup>3a</sup>, SO<sub>2</sub>NR<sup>3</sup>R<sup>3a</sup>, NR<sup>3</sup>SO<sub>2</sub>-C<sub>1-4</sub> alkyl, NR<sup>3</sup>SO<sub>2</sub>-phenyl, S(O)<sub>2</sub>-CH<sub>3</sub>, S(O)<sub>2</sub>-phenyl, CF<sub>3</sub>, phenyl substituted with 0-2 R<sup>6</sup>, naphthyl substituted with 0-2 R<sup>6</sup>, and benzyl substituted with 0-2 R<sup>6</sup>; and,

$R^6$ , at each occurrence, is selected from H, OH, OR<sup>2</sup>, F, Cl, CH<sub>3</sub>, CH<sub>2</sub>CH<sub>3</sub>, NR<sup>2</sup>R<sup>2a</sup>, CH<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>, C(O)R<sup>2b</sup>, CH<sub>2</sub>C(O)R<sup>2b</sup>, NR<sup>2</sup>C(O)R<sup>2b</sup>, and SO<sub>2</sub>NR<sup>2</sup>R<sup>2a</sup>.

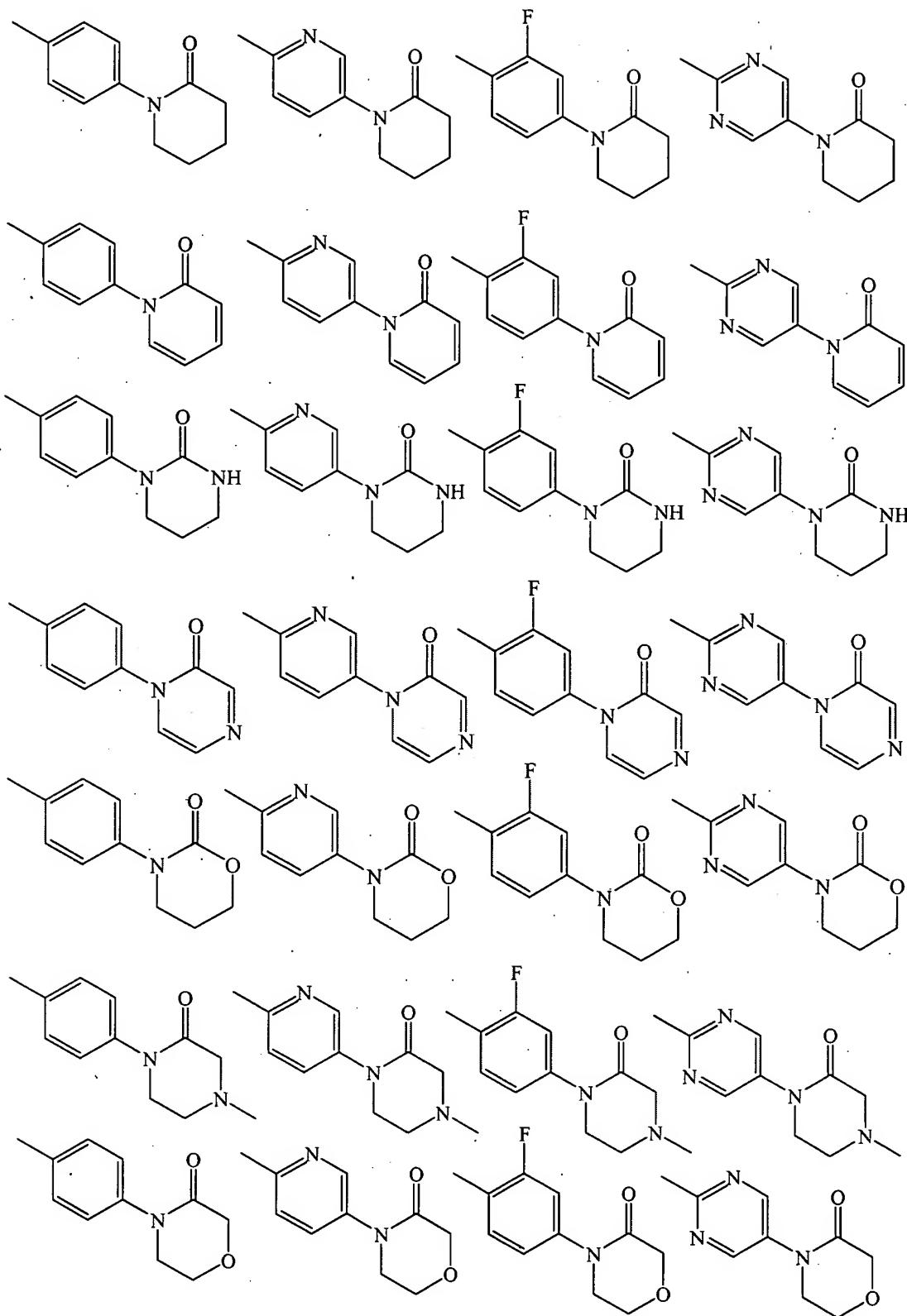
6. (Original) A compound according to Claim 5, wherein:

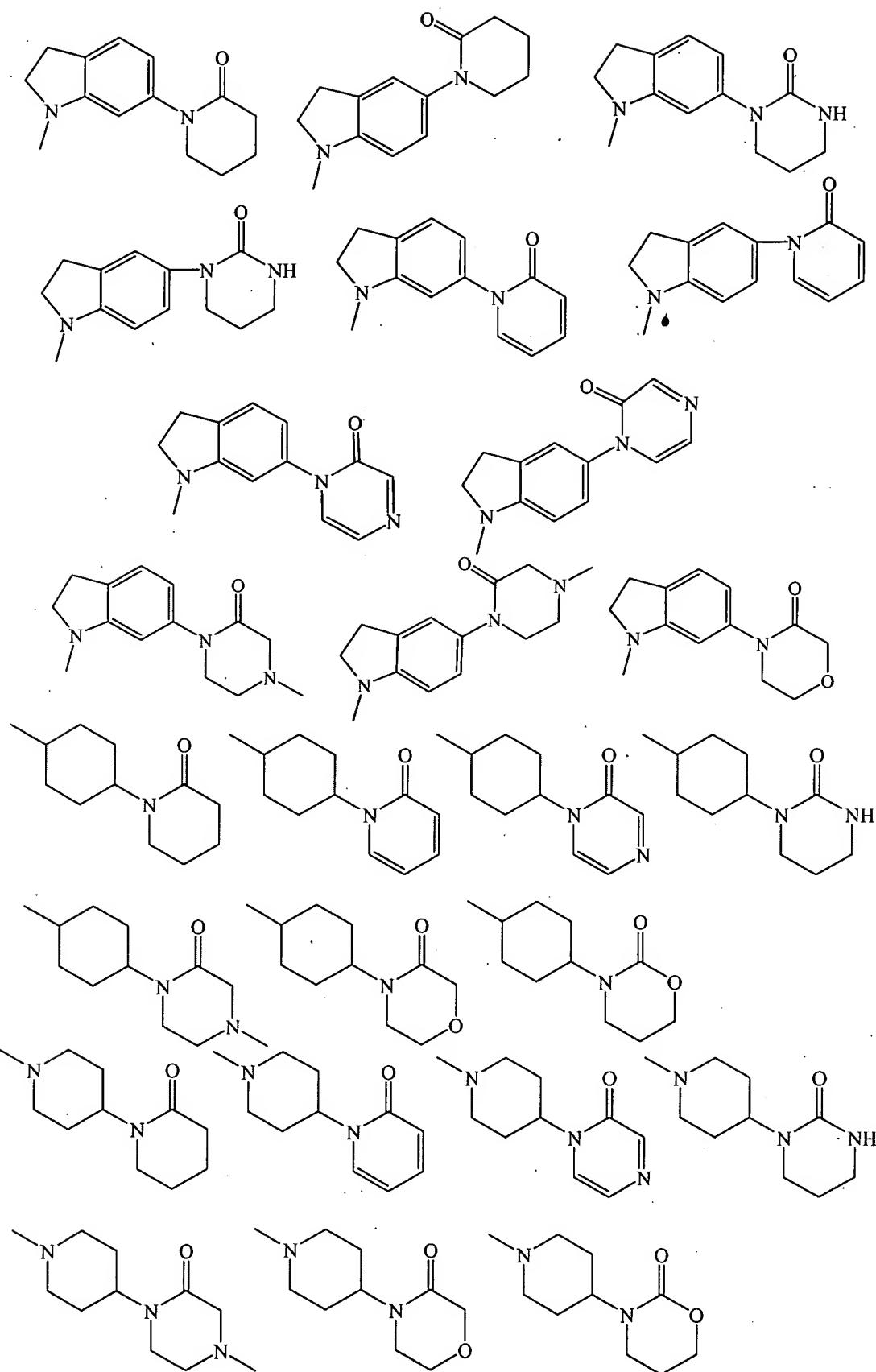
G is selected from:



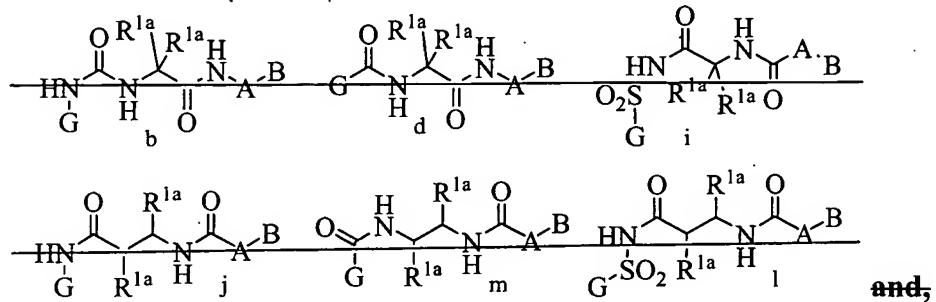


A-B is selected from:



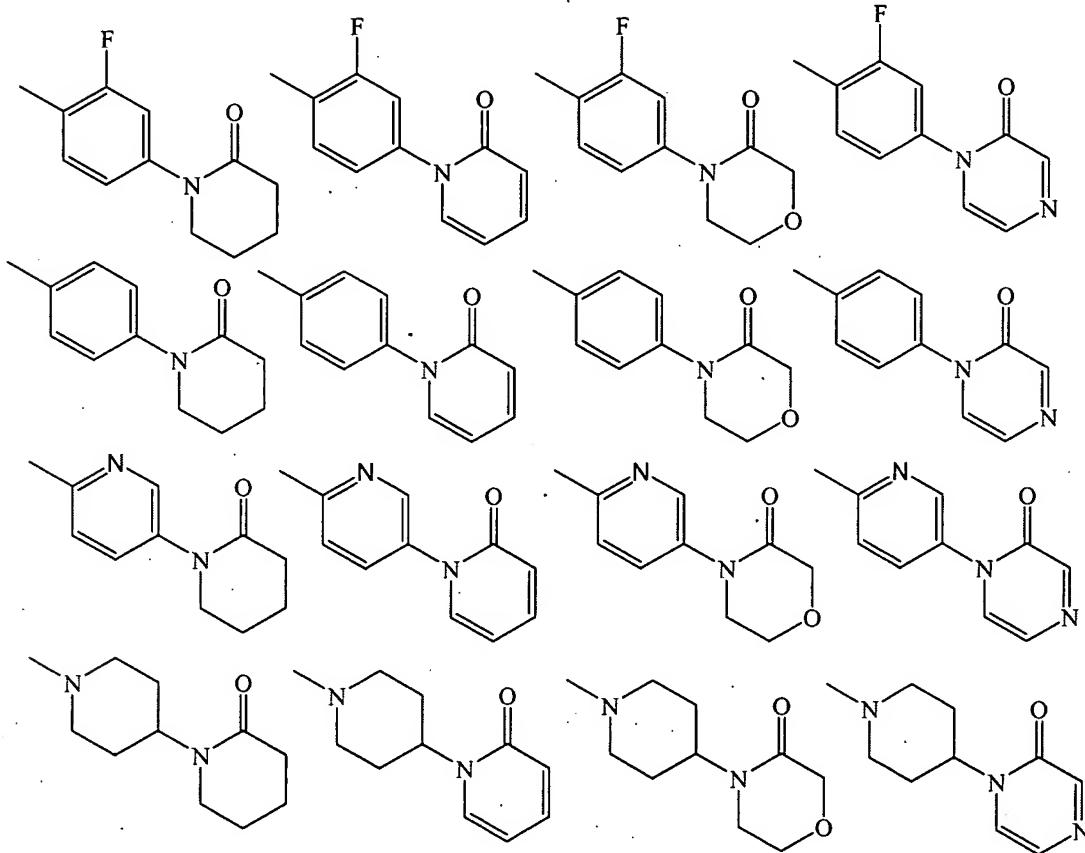


7. (Currently Amended) A compound according to Claim 6, wherein ~~the compound is selected from:~~



and,

A-B is selected from:



8. (Currently Amended) A compound ~~according to Claim 1, wherein the compound is~~ selected from the group:

2-(5-Chloro-thiophene-2-sulfonylamino)-*N*-[4-(2-oxo-2*H*-pyridin-1-yl)-phenyl]-2-phenyl-acetamide;

2-(6-Chloro-naphthalene-2-sulfonylamino)-*N*-[4-(2-oxo-2*H*-pyridin-1-yl)-phenyl]-2-phenyl-acetamide;

5-Chloro-thiophene-2-carboxylic acid {[4-(2-oxo-2*H*-pyridin-1-yl)-phenylcarbamoyl]-phenyl-methyl}-amide;

5-Chloro-1*H*-indole-2-carboxylic acid {[4-(2-oxo-2*H*-pyridin-1-yl)-phenylcarbamoyl]-phenyl-methyl}-amide;

3-Chloro-1*H*-indole-6-carboxylic acid {[4-(2-oxo-2*H*-pyridin-1-yl)-phenylcarbamoyl]-phenyl-methyl}-amide;

1*H*-Indole-6-carboxylic acid {[4-(2-oxo-2*H*-pyridin-1-yl)-phenylcarbamoyl]-phenyl-methyl}-amide;

2-*R*-(6-Chloro-naphthalene-2-sulfonylamino)-*N*-[4-(2-oxo-2*H*-pyridin-1-yl)-phenyl]-2-phenyl-acetamide;

2-*S*-(6-Chloro-naphthalene-2-sulfonylamino)-*N*-[4-(2-oxo-2*H*-pyridin-1-yl)-phenyl]-2-phenyl-acetamide;

2-(5-Chloro-thiophene-2-sulfonylamino)-*N*-[4-(2-oxo-2*H*-pyridin-1-yl)-phenyl]-2-phenyl-acetamide;

*N*-β-(6-chloro-naphthalene-2-sulfonylamino)-3-oxo-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

*N*-[β-(4-methoxyl-benzenesulfonylamino)-3-oxo-propyl]-4-(2-oxo-piperidin-1-yl)benzamide;

*N*-[2-(5-Chloro-pyridin-2-ylcarbamoyl)ethyl]-4-(2-oxo-2*H*-pyridin-1-yl)benzamide;

3-Chloro-1*H*-indole-6-carboxylic acid {2-[4-(2-oxo-2*H*-pyridin-1-yl)benzoylamino]ethyl} amide;

5-Chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2*H*-pyridin-1-yl)benzoylamino]ethyl} amide;

5-Chloro-1*H*-indole-2-carboxylic acid {2-[4-(2-oxo-2*H*-pyridin-1-yl)benzoylamino]ethyl} amide;

*N*-{4-[(4-Chloro-phenylcarbamoyl)-methyl]-tetrahydro-pyran-4-yl}-4-(2-oxo-2*H*-pyridin-1-yl)-benzamide;

2-[(5-Chloro-thiophene-2-carbonyl)-amino]-3-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-propionic acid methyl ester;  
or a pharmaceutically acceptable salt form thereof.

9. (Currently Amended) A compound according to Claim 1, wherein the compound is selected from: **Examples 19-454 of Table 1.**

**3-chloro-N-(2-{[4-(2-oxopiperidin-1-yl)benzoyl]amino}ethyl)-1H-indole-6-carboxamide;**

**5-chloro-N-(2-{[4-(2-oxopiperidin-1-yl)benzoyl]amino}ethyl)thiophene-2-carboxamide;**

**5-methoxy-N-(2-{[4-(2-oxopiperidin-1-yl)benzoyl]amino}ethyl)thiophene-2-carboxamide;**

**3-chloro-5-methoxy-N-(2-{[4-(2-oxopiperidin-1-yl)benzoyl]amino}ethyl)thiophene-2-carboxamide;**

**3,5-dichloro-N-(2-{[4-(2-oxopiperidin-1-yl)benzoyl]amino}ethyl)thiophene-2-carboxamide;**

**5-chloro-N-(2-{[4-(2-oxopiperidin-1-yl)benzoyl]amino}propyl)thiophene-2-carboxamide;**

**5-chloro-N-(3-methyl-2-{[4-(2-oxopiperidin-1-yl)benzoyl]amino}butyl)thiophene-2-carboxamide;**

**5-chloro-N-(2-{[4-(2-oxopiperidin-1-yl)benzoyl]amino}-2-phenylethyl)thiophene-2-carboxamide;**

**5-chloro-N-(2-{[4-(2-oxopiperidin-1-yl)benzoyl]amino}-3-phenylpropyl)thiophene-2-carboxamide;**

**methyl 3-{[(5-chloro-2-thienyl)carbonyl]amino}-N-[4-(2-oxopiperidin-1-yl)benzoyl]alaninate;**

**3-{[(5-chloro-2-thienyl)carbonyl]amino}-N-[4-(2-oxopiperidin-1-yl)benzoyl]alanine;**

**methyl 4-{[(5-chloro-2-thienyl)carbonyl]amino}-3-{[4-(2-oxopiperidin-1-yl)benzoyl]amino}butanoate;**

4-{{(5-chloro-2-thienyl)carbonyl}amino}-3-{{4-(2-oxopiperidin-1-yl)benzoyl}amino}butanoic acid;  
N-{{3-[(4-chlorophenyl)amino]-1-methyl-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-{{3-[(4-chlorophenyl)amino]-3-oxo-1-phenylpropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-{{1-benzyl-3-[(4-chlorophenyl)amino]-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;  
5-[(4-chlorophenyl)amino]-5-oxo-3-{{4-(2-oxopyridin-1(2H)-yl)benzoyl}amino}pentanoic acid;  
N<sup>4</sup>-{(4-chlorophenyl)-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]asparagine;  
6-[(4-chlorophenyl)amino]-6-oxo-4-{{4-(2-oxopyridin-1(2H)-yl)benzoyl}amino}hexanoic acid;  
N<sup>4</sup>-{(4-chlorophenyl)-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]aspartamide;  
N<sup>4</sup>-{(4-chlorophenyl)-N<sup>1</sup>,N<sup>1</sup>-dimethyl-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]aspartamide;  
N-{{3-[(4-chlorophenyl)amino]-1-[(dimethylamino)methyl]-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-{{3-[(4-chlorophenyl)amino]-1-[(methylamino)methyl]-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-{{1-(aminomethyl)-3-[(4-chlorophenyl)amino]-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-{{1-[(acetylamino)methyl]-3-[(4-chlorophenyl)amino]-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-{{3-[(4-chlorophenyl)amino]-1-{{[(methylamino)carbonyl]amino}methyl}-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-{{3-[(4-chlorophenyl)amino]-1-{{[(methylsulfonyl)amino]methyl}-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-{{3-[(4-chlorophenyl)amino]-1-(hydroxymethyl)-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-[(2-methoxyethoxy)methyl]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-(methoxymethyl)-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-(3-[(4-chlorophenyl)amino]-1-{[2-(dimethylamino)ethoxy]methyl}-3-oxopropyl)-4-(2-oxopyridin-1(2H)-yl)benzamide;

N<sup>4</sup>-(4-chlorophenyl)-N<sup>1</sup>-[2-(dimethylamino)ethyl]-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]aspartamide;

N<sup>4</sup>-(4-chlorophenyl)-N<sup>1</sup>-(2-morpholin-4-ylethyl)-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]aspartamide;

N<sup>4</sup>-(4-chlorophenyl)-N<sup>1</sup>-[2-(1,1-dioxidothiomorpholin-4-yl)ethyl]-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]aspartamide;

N<sup>4</sup>-(4-chlorophenyl)-N<sup>1</sup>-[2-(4-methylpiperazin-1-yl)ethyl]-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]aspartamide;

N-[3-[(4-chlorophenyl)amino]-1-({[2-(4-methylpiperazin-1-yl)ethyl]amino}methyl)-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-({methyl}[2-(4-methylpiperazin-1-yl)ethyl]amino{methyl})-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-(3-[(4-chlorophenyl)amino]-1-{[methyl(2-morpholin-4-ylethyl)amino]methyl}-3-oxopropyl)-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-(3-[(4-chlorophenyl)amino]-1-{[[2-(1,1-dioxidothiomorpholin-4-yl)ethyl](methyl)amino]methyl}-3-oxopropyl)-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-(3-[(4-chlorophenyl)amino]-1-{[2-(1,1-dioxidothiomorpholin-4-yl)ethoxy]methyl}-3-oxopropyl)-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-[(2-morpholin-4-ylethoxy)methyl]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-(3-[(4-chlorophenyl)amino]-1-{[2-(4-methylpiperazin-1-yl)ethoxy]methyl}-3-oxopropyl)-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-3-oxo-1-(pyrrolidin-1-ylcarbonyl)propyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-3-oxo-1-(piperidin-1-ylcarbonyl)propyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-(morpholin-4-ylcarbonyl)-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-{3-[(4-chlorophenyl)amino]-1-[(4-methylpiperazin-1-yl)carbonyl]-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-{3-[(4-chlorophenyl)amino]-1-[(1,1-dioxidothiomorpholin-4-yl)carbonyl]-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-(morpholin-4-ylmethyl)-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-3-oxo-1-(pyrrolidin-1-ylmethyl)propyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-{3-[(4-chlorophenyl)amino]-3-oxo-1-[(2-oxopyrrolidin-1-yl)methyl]propyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-{3-[(5-chloropyridin-2-yl)amino]-1-methyl-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-{3-[(5-chloropyridin-2-yl)amino]-3-oxo-1-phenylpropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-{1-benzyl-3-[(5-chloropyridin-2-yl)amino]-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;

5-[(5-chloropyridin-2-yl)amino]-5-oxo-3-{[4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}pentanoic acid;

N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]asparagine;

6-[(5-chloropyridin-2-yl)amino]-6-oxo-4-{[4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}hexanoic acid;

N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]aspartamide;

N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>1</sup>,N<sup>1</sup>-dimethyl-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]aspartamide;

N-{3-[(5-chloropyridin-2-yl)amino]-1-[(dimethylamino)methyl]-3-oxopropyl}-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-[3-[(5-chloropyridin-2-yl)amino]-1-[(methylamino)methyl]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[1-(aminomethyl)-3-[(5-chloropyridin-2-yl)amino]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[1-[(acetylamino)methyl]-3-[(5-chloropyridin-2-yl)amino]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-{[(methylamino)carbonyl]amino}methyl]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-(3-[(5-chloropyridin-2-yl)amino]-1-{[(methylsulfonyl)amino]methyl}-3-oxopropyl)-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-(hydroxymethyl)-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-(2-methoxyethoxy)methyl]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-(methoxymethyl)-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-(3-[(5-chloropyridin-2-yl)amino]-1-{[2-(dimethylamino)ethoxy]methyl}-3-oxopropyl)-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>1</sup>-[2-(dimethylamino)ethyl]-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]aspartamide;  
N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>1</sup>-[2-morpholin-4-ylethyl]-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]aspartamide;  
N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>1</sup>-[2-(1,1-dioxidothiomorpholin-4-yl)ethyl]-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]aspartamide;  
N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>1</sup>-[2-(4-methylpiperazin-1-yl)ethyl]-N<sup>2</sup>-[4-(2-oxopyridin-1(2H)-yl)benzoyl]aspartamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-({[2-(4-methylpiperazin-1-yl)ethyl]amino}methyl)-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-({methyl}[2-(4-methylpiperazin-1-yl)ethyl]amino)methyl)-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-[3-[(5-chloropyridin-2-yl)amino]-1-[[methyl(2-morpholin-4-ylethyl)amino]methyl]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-[[2-(1,1-dioxidothiomorpholin-4-yl)ethyl](methyl)amino]methyl]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-[[2-(1,1-dioxidothiomorpholin-4-yl)ethoxy]methyl]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-[(2-morpholin-4-ylethoxy)methyl]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-[[2-(4-methylpiperazin-1-yl)ethoxy]methyl]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-3-oxo-1-(pyrrolidin-1-ylcarbonyl)propyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-3-oxo-1-(piperidin-1-ylcarbonyl)propyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-(morpholin-4-ylcarbonyl)-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-[(4-methylpiperazin-1-yl)carbonyl]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-[(1,1-dioxidothiomorpholin-4-yl)carbonyl]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-(morpholin-4-ylmethyl)-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-3-oxo-1-(pyrrolidin-1-ylmethyl)propyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-3-oxo-1-[(2-oxopyrrolidin-1-yl)methyl]propyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(4-chlorophenyl)amino]-2-methyl-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;  
N-[3-[(4-chlorophenyl)amino]-3-oxo-2-phenylpropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-[2-benzyl-3-[(4-chlorophenyl)amino]-3-oxopropyl]-4-(2-oxopyridin-1(2H)-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-methyl-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-3-oxo-1-phenylpropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[1-benzyl-3-[(4-chlorophenyl)amino]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

5-[(4-chlorophenyl)amino]-5-oxo-3-{|4-(2-oxopiperidin-1-yl)benzoyl]amino}pentanoic acid;

N<sup>4</sup>-(4-chlorophenyl)-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]asparagine;

6-[(4-chlorophenyl)amino]-6-oxo-4-{|4-(2-oxopiperidin-1-yl)benzoyl]amino}hexanoic acid;

N<sup>4</sup>-(4-chlorophenyl)-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]aspartamide;

N<sup>4</sup>-(4-chlorophenyl)-N<sup>1</sup>,N<sup>1</sup>-dimethyl-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]aspartamide;

N-[3-[(4-chlorophenyl)amino]-1-[(dimethylamino)methyl]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-[(methylamino)methyl]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[1-(aminomethyl)-3-[(4-chlorophenyl)amino]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[1-[(acetylamino)methyl]-3-[(4-chlorophenyl)amino]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-{|[(methylamino)carbonyl]amino|methyl}-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-(3-[(4-chlorophenyl)amino]-1-{|[(methylsulfonyl)amino|methyl}-3-oxopropyl)-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-(hydroxymethyl)-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-[(2-methoxyethoxy)methyl]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-(methoxymethyl)-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-(3-[(4-chlorophenyl)amino]-1-[[2-(dimethylamino)ethoxy]methyl]-3-oxopropyl)-4-(2-oxopiperidin-1-yl)benzamide;

N<sup>4</sup>-(4-chlorophenyl)-N<sup>1</sup>-[2-(dimethylamino)ethyl]-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]aspartamide;

N<sup>4</sup>-(4-chlorophenyl)-N<sup>1</sup>-(2-morpholin-4-ylethyl)-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]aspartamide;

N<sup>4</sup>-(4-chlorophenyl)-N<sup>1</sup>-[2-(1,1-dioxidothiomorpholin-4-yl)ethyl]-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]aspartamide;

N<sup>4</sup>-(4-chlorophenyl)-N<sup>1</sup>-[2-(4-methylpiperazin-1-yl)ethyl]-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]aspartamide;

N-[3-[(4-chlorophenyl)amino]-1-{{[2-(4-methylpiperazin-1-yl)ethyl]amino}methyl}-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-{{methyl}[2-(4-methylpiperazin-1-yl)ethyl]amino}methyl]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-(3-[(4-chlorophenyl)amino]-1-{{methyl}(2-morpholin-4-ylethyl)amino}methyl}-3-oxopropyl)-4-(2-oxopiperidin-1-yl)benzamide;

N-(3-[(4-chlorophenyl)amino]-1-{{[2-(1,1-dioxidothiomorpholin-4-yl)ethyl](methyl)amino}methyl}-3-oxopropyl)-4-(2-oxopiperidin-1-yl)benzamide;

N-(3-[(4-chlorophenyl)amino]-1-{{[2-(1,1-dioxidothiomorpholin-4-yl)ethoxy]methyl}-3-oxopropyl)-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-1-[(2-morpholin-4-ylethoxy)methyl]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-(3-[(4-chlorophenyl)amino]-1-{{[2-(4-methylpiperazin-1-yl)ethoxy]methyl}-3-oxopropyl)-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-3-oxo-1-(pyrrolidin-1-ylcarbonyl)propyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(4-chlorophenyl)amino]-3-oxo-1-(piperidin-1-ylcarbonyl)propyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(4-chlorophenyl)amino]-1-(morpholin-4-ylcarbonyl)-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-{3-[(4-chlorophenyl)amino]-1-[(4-methylpiperazin-1-yl)carbonyl]-3-oxopropyl}-4-(2-oxopiperidin-1-yl)benzamide;  
N-{3-[(4-chlorophenyl)amino]-1-[(1,1-dioxidothiomorpholin-4-yl)carbonyl]-3-oxopropyl}-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(4-chlorophenyl)amino]-1-(morpholin-4-ylmethyl)-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(4-chlorophenyl)amino]-3-oxo-1-(pyrrolidin-1-ylmethyl)propyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-{3-[(4-chlorophenyl)amino]-3-oxo-1-[(2-oxopyrrolidin-1-yl)methyl]propyl}-4-(2-oxopiperidin-1-yl)benzamide;  
N-{3-[(5-chloropyridin-2-yl)amino]-1-methyl-3-oxopropyl}-4-(2-oxopiperidin-1-yl)benzamide;  
N-{3-[(5-chloropyridin-2-yl)amino]-3-oxo-1-phenylpropyl}-4-(2-oxopiperidin-1-yl)benzamide;  
N-{1-benzyl-3-[(5-chloropyridin-2-yl)amino]-3-oxopropyl}-4-(2-oxopiperidin-1-yl)benzamide;  
5-[(5-chloropyridin-2-yl)amino]-5-oxo-3-{[4-(2-oxopiperidin-1-yl)benzoyl]amino}pentanoic acid;  
N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]asparagine;  
6-[(5-chloropyridin-2-yl)amino]-6-oxo-4-{[4-(2-oxopiperidin-1-yl)benzoyl]amino}hexanoic acid;  
N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]aspartamide;  
N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>1</sup>,N<sup>1</sup>-dimethyl-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]aspartamide;  
N-{3-[(5-chloropyridin-2-yl)amino]-1-[(dimethylamino)methyl]-3-oxopropyl}-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(5-chloropyridin-2-yl)amino]-1-[(methylamino)methyl]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[1-(aminomethyl)-3-[(5-chloropyridin-2-yl)amino]-3-oxopropyl]-4-(2-oxopiperidin-1-yl benzamide;

N-[1-[(acetylamino)methyl]-3-[(5-chloropyridin-2-yl)amino]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(5-chloropyridin-2-yl)amino]-1-{[(methylamino)carbonyl]amino}methyl]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-(3-[(5-chloropyridin-2-yl)amino]-1-{[(methylsulfonyl)amino]methyl}-3-oxopropyl)-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(5-chloropyridin-2-yl)amino]-1-(hydroxymethyl)-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(5-chloropyridin-2-yl)amino]-1-[(2-methoxyethoxy)methyl]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(5-chloropyridin-2-yl)amino]-1-(methoxymethyl)-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-(3-[(5-chloropyridin-2-yl)amino]-1-{[2-(dimethylamino)ethoxy]methyl}-3-oxopropyl)-4-(2-oxopiperidin-1-yl)benzamide;

N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>1</sup>-[2-(dimethylamino)ethyl]-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]aspartamide;

N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>1</sup>-[2-morpholin-4-ylethyl]-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]aspartamide;

N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>1</sup>-[2-(1,1-dioxidothiomorpholin-4-yl)ethyl]-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]aspartamide;

N<sup>4</sup>-(5-chloropyridin-2-yl)-N<sup>1</sup>-[2-(4-methylpiperazin-1-yl)ethyl]-N<sup>2</sup>-[4-(2-oxopiperidin-1-yl)benzoyl]aspartamide;

N-[3-[(5-chloropyridin-2-yl)amino]-1-({[2-(4-methylpiperazin-1-yl)ethyl]amino}methyl)-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-[3-[(5-chloropyridin-2-yl)amino]-1-({methyl}[2-(4-methylpiperazin-1-yl)ethyl]amino}methyl)-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-(3-[(5-chloropyridin-2-yl)amino]-1-{[methyl(2-morpholin-4-ylethyl)amino]methyl}-3-oxopropyl)-4-(2-oxopiperidin-1-yl)benzamide;  
N-(3-[(5-chloropyridin-2-yl)amino]-1-{[2-(1,1-dioxidothiomorpholin-4-yl)ethyl](methyl)amino]methyl}-3-oxopropyl)-4-(2-oxopiperidin-1-yl)benzamide;  
N-(3-[(5-chloropyridin-2-yl)amino]-1-{[2-(1,1-dioxidothiomorpholin-4-yl)ethoxy]methyl}-3-oxopropyl)-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-[(2-morpholin-4-ylethoxy)methyl]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-(3-[(5-chloropyridin-2-yl)amino]-1-{[2-(4-methylpiperazin-1-yl)ethoxy]methyl}-3-oxopropyl)-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-3-oxo-1-(pyrrolidin-1-ylcarbonyl)propyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-3-oxo-1-(piperidin-1-ylcarbonyl)propyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-(morpholin-4-ylcarbonyl)-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-[(4-methylpiperazin-1-yl)carbonyl]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-[(1,1-dioxidothiomorpholin-4-yl)carbonyl]-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-1-(morpholin-4-ylmethyl)-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-3-oxo-1-(pyrrolidin-1-ylmethyl)propyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(5-chloropyridin-2-yl)amino]-3-oxo-1-(2-oxopyrrolidin-1-yl)methyl]propyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(4-chlorophenyl)amino]-2-methyl-3-oxopropyl]-4-(2-oxopiperidin-1-yl)benzamide;  
N-[3-[(4-chlorophenyl)amino]-3-oxo-2-phenylpropyl]-4-(2-oxopiperidin-1-yl)benzamide;

N-{2-benzyl-3-[(4-chlorophenyl)amino]-3-oxopropyl}-4-(2-oxopiperidin-1-yl)benzamide;  
5-chloro-N-(2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl|amino}ethyl)thiophene-2-carboxamide;  
5-chloro-N-(2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl|amino}propyl)thiophene-2-carboxamide;  
5-chloro-N-(2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl|amino}-2-phenylethyl)thiophene-2-carboxamide;  
5-chloro-N-(2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl|amino}-3-phenylpropyl)thiophene-2-carboxamide;  
3-{|(5-chloro-2-thienyl)carbonyl|amino}-N-[4-(2-oxopyridin-1(2H)-yl)benzoyl]alanine;  
4-{|(5-chloro-2-thienyl)carbonyl|amino}-3-{|4-(2-oxopyridin-1(2H)-yl)benzoyl|amino}butanoic acid;  
5-{|(5-chloro-2-thienyl)carbonyl|amino}-4-{|4-(2-oxopyridin-1(2H)-yl)benzoyl|amino}pentanoic acid;  
N-(3-amino-2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl|amino}propyl)-5-chlorothiophene-2-carboxamide;  
5-chloro-N-(3-(methylamino)-2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl|amino}propyl)thiophene-2-carboxamide;  
5-chloro-N-(3-(dimethylamino)-2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl|amino}propyl)thiophene-2-carboxamide;  
5-chloro-N-(3-|[2-(dimethylamino)ethyl](methyl)amino]-2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl|amino}propyl)thiophene-2-carboxamide;  
5-chloro-N-(3-[methyl(2-morpholin-4-ylethyl)amino]-2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl|amino}propyl)thiophene-2-carboxamide;  
5-chloro-N-(3-|[2-(1,1-dioxidothiomorpholin-4-yl)ethyl](methyl)amino]-2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl|amino}propyl)thiophene-2-carboxamide;  
5-chloro-N-(3-|[2-(1,1-dioxidothiomorpholin-4-yl)ethoxy]-2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl|amino}propyl)thiophene-2-carboxamide;

5-chloro-N-(3-(2-morpholin-4-ylethoxy)-2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}propyl)thiophene-2-carboxamide;  
5-chloro-N-(3-[2-(2-oxopiperidin-1-yl)ethoxy]-2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}propyl)thiophene-2-carboxamide;  
5-chloro-N-(1-methyl-2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}ethyl)thiophene-2-carboxamide;  
5-chloro-N-(2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}-1-phenylethyl)thiophene-2-carboxamide;  
N-(1-benzyl-2-{|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}ethyl)-5-chlorothiophene-2-carboxamide;  
3-{|[(5-chloro-2-thienyl)carbonyl]amino}-4-{|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}butanoic acid;  
N-[(5-chloro-2-thienyl)carbonyl]-3-{|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}alanine;  
N-[2-amino-1-({|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]-5-chlorothiophene-2-carboxamide;  
5-chloro-N-[2-(methylamino)-1-({|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-(dimethylamino)-1-({|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;  
N-[2-[acetyl(methyl)amino]-1-({|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]-5-chlorothiophene-2-carboxamide;  
5-chloro-N-[2-[methyl(methylsulfonyl)amino]-1-({|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-hydroxy-1-({|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-methoxy-1-({|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-[2-(dimethylamino)ethoxy]-1-({|4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;

5-chloro-N-[2-(2-morpholin-4-ylethoxy)-1-({[4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-[2-(1,1-dioxidothiomorpholin-4-yl)ethoxy]-1-({[4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-[2-(4-methylpiperazin-1-yl)ethoxy]-1-({[4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-{methyl}[2-(4-methylpiperazin-1-yl)ethyl]amino}-1-({[4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-[methyl(2-morpholin-4-ylethyl)amino]-1-({[4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-[2-(1,1-dioxidothiomorpholin-4-yl)ethyl](methyl)amino]-1-({[4-(2-oxopyridin-1(2H)-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;  
5-chloro-N-(2-{{[4-(2-oxopiperidin-1-yl)benzoyl]amino}ethyl}thiophene-2-carboxamide;  
5-chloro-N-(2-{{[4-(2-oxopiperidin-1-yl)benzoyl]amino}propyl}thiophene-2-carboxamide;  
5-chloro-N-(2-{{[4-(2-oxopiperidin-1-yl)benzoyl]amino}-2-phenylethyl}thiophene-2-carboxamide;  
5-chloro-N-(2-{{[4-(2-oxopiperidin-1-yl)benzoyl]amino}-3-phenylpropyl}thiophene-2-carboxamide;  
3-{{[(5-chloro-2-thienyl)carbonyl]amino}-N-[4-(2-oxopiperidin-1-yl)benzoyl]alanine};  
4-{{[(5-chloro-2-thienyl)carbonyl]amino}-3-{{[4-(2-oxopiperidin-1-yl)benzoyl]amino}butanoic acid};  
5-{{[(5-chloro-2-thienyl)carbonyl]amino}-4-{{[4-(2-oxopiperidin-1-yl)benzoyl]amino}pentanoic acid};  
N-(3-amino-2-{{[4-(2-oxopiperidin-1-yl)benzoyl]amino}propyl}-5-chlorothiophene-2-carboxamide;  
5-chloro-N-(3-(methylamino)-2-{{[4-(2-oxopiperidin-1-yl)benzoyl]amino}propyl}thiophene-2-carboxamide;  
5-chloro-N-(3-(dimethylamino)-2-{{[4-(2-oxopiperidin-1-yl)benzoyl]amino}propyl}thiophene-2-carboxamide;

5-chloro-N-(3-[2-(dimethylamino)ethyl](methyl)amino]-2-{|4-(2-oxopiperidin-1-yl)benzoyl]amino}propyl)thiophene-2-carboxamide;

5-chloro-N-(3-[methyl(2-morpholin-4-ylethyl)amino]-2-{|4-(2-oxopiperidin-1-yl)benzoyl]amino}propyl)thiophene-2-carboxamide;

5-chloro-N-(3-[2-(1,1-dioxidothiomorpholin-4-yl)ethyl](methyl)amino]-2-{|4-(2-oxopiperidin-1-yl)benzoyl]amino}propyl)thiophene-2-carboxamide;

5-chloro-N-(3-[2-(1,1-dioxidothiomorpholin-4-yl)ethoxy]-2-{|4-(2-oxopiperidin-1-yl)benzoyl]amino}propyl)thiophene-2-carboxamide;

5-chloro-N-(3-(2-morpholin-4-ylethoxy)-2-{|4-(2-oxopiperidin-1-yl)benzoyl]amino}propyl)thiophene-2-carboxamide;

5-chloro-N-(3-[2-(2-oxopiperidin-1-yl)ethoxy]-2-{|4-(2-oxopiperidin-1-yl)benzoyl]amino}propyl)thiophene-2-carboxamide;

5-chloro-N-(1-methyl-2-{|4-(2-oxopiperidin-1-yl)benzoyl]amino}ethyl)thiophene-2-carboxamide;

5-chloro-N-(2-{|4-(2-oxopiperidin-1-yl)benzoyl]amino}-1-phenylethyl)thiophene-2-carboxamide;

N-(1-benzyl-2-{|4-(2-oxopiperidin-1-yl)benzoyl]amino}ethyl)-5-chlorothiophene-2-carboxamide;

3-{|[(5-chloro-2-thienyl)carbonyl]amino}-4-{|4-(2-oxopiperidin-1-yl)benzoyl]amino}butanoic acid;

N-[(5-chloro-2-thienyl)carbonyl]-3-{|4-(2-oxopiperidin-1-yl)benzoyl]amino}alanine;

N-[2-amino-1-({|4-(2-oxopiperidin-1-yl)benzoyl]amino}methyl)ethyl]-5-chlorothiophene-2-carboxamide;

5-chloro-N-[2-(methylamino)-1-({|4-(2-oxopiperidin-1-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;

5-chloro-N-[2-(dimethylamino)-1-({|4-(2-oxopiperidin-1-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;

N-[2-[acetyl(methyl)amino]-1-({|4-(2-oxopiperidin-1-yl)benzoyl]amino}methyl)ethyl]-5-chlorothiophene-2-carboxamide;

5-chloro-N-[2-[methyl(methylsulfonyl)amino]-1-({|4-(2-oxopiperidin-1-yl)benzoyl]amino}methyl)ethyl]thiophene-2-carboxamide;

5-chloro-N-[2-hydroxy-1-([4-(2-oxopiperidin-1-yl)benzoyl]amino)methyl]ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-methoxy-1-([4-(2-oxopiperidin-1-yl)benzoyl]amino)methyl]ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-[2-(dimethylamino)ethoxy]-1-([4-(2-oxopiperidin-1-yl)benzoyl]amino)methyl]ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-(2-morpholin-4-ylethoxy)-1-([4-(2-oxopiperidin-1-yl)benzoyl]amino)methyl]ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-[2-(1,1-dioxidothiomorpholin-4-yl)ethoxy]-1-([4-(2-oxopiperidin-1-yl)benzoyl]amino)methyl]ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-[2-(4-methylpiperazin-1-yl)ethoxy]-1-([4-(2-oxopiperidin-1-yl)benzoyl]amino)methyl]ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-{methyl}[2-(4-methylpiperazin-1-yl)ethyl]amino]-1-([4-(2-oxopiperidin-1-yl)benzoyl]amino)methyl]ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-[methyl(2-morpholin-4-ylethyl)amino]-1-([4-(2-oxopiperidin-1-yl)benzoyl]amino)methyl]ethyl]thiophene-2-carboxamide;  
5-chloro-N-[2-[2-(1,1-dioxidothiomorpholin-4-yl)ethyl](methyl)amino]-1-([4-(2-oxopiperidin-1-yl)benzoyl]amino)methyl]ethyl]thiophene-2-carboxamide;  
3-chloro-N-(2-oxo-2-[(4-(2-oxopiperidin-1-yl)phenyl)amino]-1-phenylethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(2-fluorophenyl)-2-oxo-2-[(4-(2-oxopiperidin-1-yl)phenyl)amino]ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(2-oxo-2-[(4-(2-oxopyridin-1(2H)-yl)phenyl)amino]-1-phenylethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(2-chlorophenyl)-2-oxo-2-[(4-(2-oxopyridin-1(2H)-yl)phenyl)amino]ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(2-fluorophenyl)-2-oxo-2-[(4-(2-oxopyridin-1(2H)-yl)phenyl)amino]ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(2-methoxyphenyl)-2-oxo-2-[(4-(2-oxopyridin-1(2H)-yl)phenyl)amino]ethyl)-1H-indole-6-carboxamide;

N-(1-[3-(aminocarbonyl)phenyl]-2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}ethyl)-3-chloro-1H-indole-6-carboxamide;  
3-chloro-N-(1-[3-(methylsulfonyl)phenyl]-2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(2-methylphenyl)-2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}-1-pyridin-2-ylethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}-1-pyridin-3-ylethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}-1-pyridin-4-ylethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(2-cyanophenyl)-2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-[2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}-1-(3-thienyl)ethyl]-1H-indole-6-carboxamide;  
3-chloro-N-[2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}-1-(2-thienyl)ethyl]-1H-indole-6-carboxamide;  
3-chloro-N-[2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}-1-(4-thienyl)ethyl]-1H-indole-6-carboxamide;  
3-chloro-N-[2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}-1-(1,3-thiazol-4-yl)ethyl]-1H-indole-6-carboxamide;  
3-chloro-N-[2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}-1-(1,3-thiazol-5-yl)ethyl]-1H-indole-6-carboxamide;  
3-chloro-N-[2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}-1-(1,3-thiazol-2-yl)ethyl]-1H-indole-6-carboxamide;  
3-chloro-N-(1-(1-methyl-1H-pyrazol-4-yl)-2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(2-naphthyl)-2-oxo-2-{\[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;

3-chloro-N-(1-(1-naphthyl)-2-oxo-2-{|4-(2-oxopyridin-1(2H)-yl)phenyl]amino}ethyl}-1H-indole-6-carboxamide;  
N-(1-(1-benzothien-2-yl)-2-oxo-2-{|4-(2-oxopyridin-1(2H)-yl)phenyl]amino}ethyl)-3-chloro-1H-indole-6-carboxamide;  
3-chloro-N-(2-oxo-2-{|4-(2-oxopyridin-1(2H)-yl)phenyl]amino}-1-quinolin-4-ylethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(2-oxo-2-{|4-(2-oxopiperidin-1-yl)phenyl]amino}-1-phenylethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(2-chlorophenyl)-2-oxo-2-{|4-(2-oxopiperidin-1-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(2-fluorophenyl)-2-oxo-2-{|4-(2-oxopiperidin-1-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(2-methoxyphenyl)-2-oxo-2-{|4-(2-oxopiperidin-1-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;  
N-(1-[3-(aminocarbonyl)phenyl]-2-oxo-2-{|4-(2-oxopiperidin-1-yl)phenyl]amino}ethyl)-3-chloro-1H-indole-6-carboxamide;  
3-chloro-N-(1-[3-(methylsulfonyl)phenyl]-2-oxo-2-{|4-(2-oxopiperidin-1-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(2-methylphenyl)-2-oxo-2-{|4-(2-oxopiperidin-1-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(2-oxo-2-{|4-(2-oxopiperidin-1-yl)phenyl]amino}-1-pyridin-2-ylethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(2-oxo-2-{|4-(2-oxopiperidin-1-yl)phenyl]amino}-1-pyridin-3-ylethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(2-oxo-2-{|4-(2-oxopiperidin-1-yl)phenyl]amino}-1-pyridin-4-ylethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(2-cyanophenyl)-2-oxo-2-{|4-(2-oxopiperidin-1-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-[2-oxo-2-{|4-(2-oxopiperidin-1-yl)phenyl]amino}-1-(3-thienyl)ethyl]-1H-indole-6-carboxamide;

3-chloro-N-[2-oxo-2-{\[4-(2-oxopiperidin-1-yl)phenyl]amino}-1-(2-thienyl)ethyl]-1H-indole-6-carboxamide;  
3-chloro-N-[2-oxo-2-{\[4-(2-oxopiperidin-1-yl)phenyl]amino}-1-(4-thienyl)ethyl]-1H-indole-6-carboxamide;  
3-chloro-N-[2-oxo-2-{\[4-(2-oxopiperidin-1-yl)phenyl]amino}-1-(1,3-thiazol-4-yl)ethyl]-1H-indole-6-carboxamide;  
3-chloro-N-[2-oxo-2-{\[4-(2-oxopiperidin-1-yl)phenyl]amino}-1-(1,3-thiazol-5-yl)ethyl]-1H-indole-6-carboxamide;  
3-chloro-N-[2-oxo-2-{\[4-(2-oxopiperidin-1-yl)phenyl]amino}-1-(1,3-thiazol-2-yl)ethyl]-1H-indole-6-carboxamide;  
3-chloro-N-(1-(1-methyl-1H-pyrazol-4-yl)-2-oxo-2-{\[4-(2-oxopiperidin-1-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(2-naphthyl)-2-oxo-2-{\[4-(2-oxopiperidin-1-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;  
3-chloro-N-(1-(1-naphthyl)-2-oxo-2-{\[4-(2-oxopiperidin-1-yl)phenyl]amino}ethyl)-1H-indole-6-carboxamide;  
N-(1-(1-benzothien-2-yl)-2-oxo-2-{\[4-(2-oxopiperidin-1-yl)phenyl]amino}ethyl)-3-chloro-1H-indole-6-carboxamide;  
3-chloro-N-(2-oxo-2-{\[4-(2-oxopiperidin-1-yl)phenyl]amino}-1-quinolin-4-ylethyl)-1H-indole-6-carboxamide;  
N<sup>2</sup>-[(3-chloro-1H-indol-6-yl)carbonyl]-N<sup>1</sup>-[4-(2-oxopiperidin-1-yl)phenyl]aspartamide;  
N<sup>2</sup>-[(3-chloro-1H-indol-6-yl)carbonyl]-N<sup>1</sup>-[4-(2-oxopiperidin-1-yl)phenyl]- $\alpha$ -asparagine;  
2-{\[(3-chloro-1H-indol-6-yl)carbonyl]amino}-N-[4-(2-oxopiperidin-1-yl)phenyl]malonamide;  
N<sup>2</sup>-[(3-chloro-1H-indol-6-yl)carbonyl]-N<sup>1</sup>-[4-(2-oxopiperidin-1-yl)phenyl]glutamamide;  
3-chloro-N-[3-(methylsulfonyl)-1-({[4-(2-oxopiperidin-1-yl)phenyl]amino}carbonyl)propyl]-1H-indole-6-carboxamide;  
3-chloro-N-[1-({[4-(2-oxopiperidin-1-yl)phenyl]amino}carbonyl)-3-phenylpropyl]-1H-indole-6-carboxamide;

N-[**(3-chloro-1H-indol-6-yl)carbonyl**]-N-[4-(2-oxopyridin-1(2H)-yl)phenyl]phenylalaninamide;  
N<sup>2</sup>-[**(3-chloro-1H-indol-6-yl)carbonyl**]-N<sup>1</sup>-[4-(2-oxopyridin-1(2H)-yl)phenyl]aspartamide;  
N<sup>2</sup>-[**(3-chloro-1H-indol-6-yl)carbonyl**]-N<sup>1</sup>-[4-(2-oxopyridin-1(2H)-yl)phenyl]-□-asparagine;  
2-{[**(3-chloro-1H-indol-6-yl)carbonyl**]amino}-N-[4-(2-oxopyridin-1(2H)-yl)phenyl]malonamide;  
N<sup>2</sup>-[**(3-chloro-1H-indol-6-yl)carbonyl**]-N<sup>1</sup>-[4-(2-oxopyridin-1(2H)-yl)phenyl]glutamamide;  
3-chloro-N-[3-(methylsulfonyl)-1-({[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}carbonyl)propyl]-1H-indole-6-carboxamide;  
3-chloro-N-[1-({[4-(2-oxopyridin-1(2H)-yl)phenyl]amino}carbonyl)-3-phenylpropyl]-1H-indole-6-carboxamide;  
N-[3-(6-Chloro-thieno[2,3-b]pyridine-2-sulfonylamino)-3-oxo-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;  
N-[3-(6-Chloro-thieno[2,3-b]pyridine-2-sulfonylamino)-3-oxo-propyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;  
N-[2-(6-Chloro-naphthalene-2-sulfonylamino)-2-oxo-ethyl]-4-(2-oxo-piperidin-1-yl)-benzamide;  
N-[2-(6-Chloro-naphthalene-2-sulfonylamino)-1-methyl-2-oxo-ethyl]-4-(2-oxo-piperidin-1-yl)-benzamide;  
N-[2-(6-Chloro-naphthalene-2-sulfonylamino)-1-methyl-2-oxo-ethyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;  
N-[2-(6-Chloro-naphthalene-2-sulfonylamino)-2-oxo-ethyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;  
N-[3-(6-Chloro-thieno[2,3-b]pyridine-2-sulfonylamino)-3-oxo-propyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;  
N-[3-(6-Chloro-1H-indole-2-sulfonylamino)-3-oxo-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[3-(6-Chloro-1H-indole-2-sulfonylamino)-3-oxo-propyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;

N-[3-(6-Chloro-1H-indole-2-sulfonylamino)-3-oxo-propyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[3-(4-Chloro-benzenesulfonylamino)-3-oxo-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[3-(4-Chloro-benzenesulfonylamino)-3-oxo-propyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;

N-[2-(6-Chloro-1H-indole-2-sulfonylamino)-2-oxo-ethyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[1-(6-Chloro-naphthalene-2-sulfonylaminocarbonyl)-2-methyl-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[2-(6-Chloro-naphthalene-2-sulfonylamino)-1-methyl-2-oxo-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[2-(6-Chloro-naphthalene-2-sulfonylamino)-2-oxo-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[3-(5-Chloro-pyridine-2-sulfonylamino)-3-oxo-propyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;

N-[1-(6-Chloro-1H-indole-2-sulfonylaminocarbonyl)-2-methyl-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[1-(6-Chloro-naphthalene-2-sulfonylaminocarbonyl)-2-methyl-butyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[2-(5-Chloro-thiophene-2-sulfonylamino)-1-methyl-2-oxo-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[2-(6-Chloro-thieno[2,3-b]pyridine-2-sulfonylamino)-1-methyl-2-oxo-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[2-(5-Chloro-thiophene-2-sulfonylamino)-1-methyl-2-oxo-ethyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;

N-[3-(6-Chloro-1H-indole-2-sulfonylamino)-2-methyl-3-oxo-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[3-(6-Chloro-1H-indole-2-sulfonylamino)-2-methyl-3-oxo-propyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;

N-[3-(6-Chloro-1H-indole-2-sulfonylamino)-2-methyl-3-oxo-propyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[3-(4-Chloro-benzenesulfonylamino)-2-methyl-3-oxo-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[3-(4-Chloro-benzenesulfonylamino)-2-methyl-3-oxo-propyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;

N-[3-(6-Chloro-1H-indole-2-sulfonylamino)-3-oxo-2-phenyl-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[3-(6-Chloro-1H-indole-2-sulfonylamino)-3-oxo-2-phenyl-propyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;

N-[3-(6-Chloro-1H-indole-2-sulfonylamino)-3-oxo-2-phenyl-propyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[3-(4-Chloro-benzenesulfonylamino)-3-oxo-2-phenyl-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[3-(5-Chloro-pyridine-2-sulfonylamino)-3-oxo-2-phenyl-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[2-(6-Chloro-1H-indole-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[2-(6-Chloro-naphthalene-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[2-(5-Chloro-thiophene-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[2-(6-Chloro-thieno[2,3-b]pyridine-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[2-(5-Chloro-thiophene-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;

N-[2-(4-Chloro-benzenesulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[2-(5-Chloro-pyridine-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[2-(4-Chloro-benzenesulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[2-(5-Chloro-pyridine-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[2-(4-Chloro-benzenesulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;

N-[3-(6-Chloro-benzo[b]thiophene-2-sulfonylamino)-2-methyl-3-oxo-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[3-(6-Chloro-benzo[b]thiophene-2-sulfonylamino)-2-methyl-3-oxo-propyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;

N-[3-(6-Chloro-benzo[b]thiophene-2-sulfonylamino)-2-methyl-3-oxo-propyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[2-(6-Chloro-benzo[b]thiophene-2-sulfonylamino)-1-methyl-2-oxo-ethyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[2-(6-Chloro-benzo[b]thiophene-2-sulfonylamino)-1-methyl-2-oxo-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[2-(6-Chloro-benzo[b]thiophene-2-sulfonylamino)-1-methyl-2-oxo-ethyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;

N-[3-(6-Chloro-3-methyl-benzo[b]thiophene-2-sulfonylamino)-2-methyl-3-oxo-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[3-(6-Chloro-3-methyl-benzo[b]thiophene-2-sulfonylamino)-2-methyl-3-oxo-propyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;

N-[3-(6-Chloro-3-methyl-benzo[b]thiophene-2-sulfonylamino)-2-methyl-3-oxo-propyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[2-(6-Chloro-3-methyl-benzo[b]thiophene-2-sulfonylamino)-1-methyl-2-oxo-ethyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;

N-[2-(6-Chloro-3-methyl-benzo[b]thiophene-2-sulfonylamino)-1-methyl-2-oxo-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;

N-[2-(6-Chloro-3-methyl-benzo[b]thiophene-2-sulfonylamino)-1-methyl-2-oxo-ethyl]-4-(2-oxo-piperidin-1-yl)-benzamide;  
N-[2-Methyl-3-(6-methyl-benzo[b]thiophene-2-sulfonylamino)-3-oxo-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;  
N-[2-Methyl-1-(6-methyl-benzo[b]thiophene-2-sulfonylaminocarbonyl)-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;  
N-[2-Methyl-3-(6-methyl-benzo[b]thiophene-2-sulfonylamino)-3-oxo-propyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;  
N-[2-Methyl-1-(6-methyl-benzo[b]thiophene-2-sulfonylaminocarbonyl)-propyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;  
N-[2-Methyl-3-(6-methyl-benzo[b]thiophene-2-sulfonylamino)-3-oxo-propyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;  
N-[2-Methyl-1-(6-methyl-benzo[b]thiophene-2-sulfonylaminocarbonyl)-propyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;  
N-[3-(3,6-Dimethyl-benzo[b]thiophene-2-sulfonylamino)-2-methyl-3-oxo-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;  
N-[1-(3,6-Dimethyl-benzo[b]thiophene-2-sulfonylaminocarbonyl)-2-methyl-propyl]-4-(2-oxo-piperidin-1-yl)-benzamide;  
N-[2-(3,6-Dimethyl-benzo[b]thiophene-2-sulfonylaminocarbonyl)-butyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;  
N-[1-(3,6-Dimethyl-benzo[b]thiophene-2-sulfonylaminocarbonyl)-2-methyl-propyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;  
N-[3-(6-Chloro-3-methyl-benzo[b]thiophene-2-sulfonylamino)-2-methyl-3-oxo-propyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;  
N-[2-(3,6-Dimethyl-benzo[b]thiophene-2-sulfonylamino)-1-methyl-2-oxo-ethyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;  
N-[2-(6-Chloro-benzo[b]thiophene-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-piperidin-1-yl)-benzamide;  
N-[2-(6-Methyl-benzo[b]thiophene-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-piperidin-1-yl)-benzamide;

N-[2-(6-Chloro-benzo[b]thiophene-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;  
N-[2-(6-Chloro-benzo[b]thiophene-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;  
N-[2-(6-Methyl-benzo[b]thiophene-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;  
N-[2-(6-Chloro-benzo[b]thiophene-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-piperidin-1-yl)-benzamide;  
N-[2-(3,6-Dimethyl-benzo[b]thiophene-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-piperidin-1-yl)-benzamide;  
N-[2-(6-Chloro-benzo[b]thiophene-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;  
N-[2-(3,6-Dimethyl-benzo[b]thiophene-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;  
N-[2-(3,6-Dimethyl-benzo[b]thiophene-2-sulfonylamino)-2-oxo-1-phenyl-ethyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;  
2-Oxo-3',4',5',6'-tetrahydro-2H,2'H-[1,4']bipyridinyl-1'-carboxylic acid {2-[(5-chloro-thiophene-2-carbonyl)-amino]-ethyl}-amide;  
4-(2-Oxo-2H-pyrazin-1-yl)-piperidine-1-carboxylic acid {2-[(5-chloro-thiophene-2-carbonyl)-amino]-ethyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid {[4-(2-oxo-2H-pyrazin-1-yl)-phenylcarbamoyl]-phenyl-methyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid {[4-(3-oxo-morpholin-4-yl)-phenylcarbamoyl]-phenyl-methyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid {[4-(2-oxo-[1,3]oxazinan-3-yl)-phenylcarbamoyl]-phenyl-methyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid {3-(1-methyl-1H-imidazol-2-yl)-1-[4-(2-oxo-2H-pyridin-1-yl)-phenylcarbamoyl]-propyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid {[4-(2-oxo-2H-pyridin-1-yl)-phenylcarbamoyl]-[tetrahydro-pyran-4-yl]-methyl}-amide;

5-Chloro-thiophene-2-carboxylic acid {[4-(2-oxo-2H-pyrazin-1-yl)-phenylcarbamoyl]-(tetrahydro-pyran-4-yl)-methyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyrazin-1-yl)-benzoylamino]-ethyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyrazin-1-yl)-benzoylamino]-propyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid (3-methoxy-1-[[4-(2-oxo-2H-pyrazin-1-yl)-benzoylamino]-methyl}-propyl)-amide;  
5-Chloro-thiophene-2-carboxylic acid {4-methoxy-2-[4-(2-oxo-2H-pyrazin-1-yl)-benzoylamino]-butyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid {3-(1-methyl-1H-imidazol-2-yl)-2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-propyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid {1-(1-methyl-1H-imidazol-2-ylmethyl)-2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-ethyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-2-phenyl-ethyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-1-phenyl-ethyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyrazin-1-yl)-benzoylamino]-2-phenyl-ethyl}-amide;  
5-Chloro-thiophene-2-carboxylic acid {1-methyl-2-[4-(2-oxo-2H-pyrazin-1-yl)-benzoylamino]-propyl}-amide;  
N-[2-(4-Chloro-phenylcarbamoyl)-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;  
N-[2-(4-Chloro-phenylcarbamoyl)-1-methyl-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;  
N-[2-(4-Chloro-phenylcarbamoyl)-propyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide;  
and  
N-[2-(5-Chloro-pyridin-2-ylcarbamoyl)-1-methyl-ethyl]-4-(2-oxo-2H-pyrazin-1-yl)-benzamide.

10. (Original) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 1 or a pharmaceutically acceptable salt form thereof.
11. (Withdrawn) A method for treating a thromboembolic disorder, comprising: administering to a patient in need thereof a therapeutically effective amount of a compound of Claim 1 or a pharmaceutically acceptable salt form thereof.
12. (Withdrawn) A method according to Claim 11, wherein the thromboembolic disorder is selected from arterial cardiovascular thromboembolic disorders, venous cardiovascular thromboembolic disorders, and thromboembolic disorders in the chambers of the heart.
13. (Withdrawn) A method according to Claim 11, wherein the thromboembolic disorder is selected from unstable angina, an acute coronary syndrome, first myocardial infarction, recurrent myocardial infarction, ischemic sudden death, transient ischemic attack, stroke, atherosclerosis, peripheral occlusive arterial disease, venous thrombosis, deep vein thrombosis, thrombophlebitis, arterial embolism, coronary arterial thrombosis, cerebral arterial thrombosis, cerebral embolism, kidney embolism, pulmonary embolism, and thrombosis resulting from (a) prosthetic valves or other implants, (b) indwelling catheters, (c) stents, (d) cardiopulmonary bypass, (e) hemodialysis, or (f) other procedures in which blood is exposed to an artificial surface that promotes thrombosis.
14. (Withdrawn) A method for treating a thromboembolic disorder, comprising: administering to a patient in need thereof a therapeutically effective amount of a first and second therapeutic agent, wherein the first therapeutic agent is compound of Claim 1 or a pharmaceutically acceptable salt thereof and the second therapeutic agent is at least one

agent selected from a second factor Xa inhibitor, an anti-coagulant agent, an anti-platelet agent, a thrombin inhibiting agent, a thrombolytic agent, and a fibrinolytic agent.

15. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 2 or a pharmaceutically acceptable salt form thereof.

16. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 3 or a pharmaceutically acceptable salt form thereof.

17. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 4 or a pharmaceutically acceptable salt form thereof.

18. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 5 or a pharmaceutically acceptable salt form thereof.

19. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 6 or a pharmaceutically acceptable salt form thereof.

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20. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 7 or a pharmaceutically acceptable salt form thereof.
21. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 8 or a pharmaceutically acceptable salt form thereof.
22. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 9 or a pharmaceutically acceptable salt form thereof.